

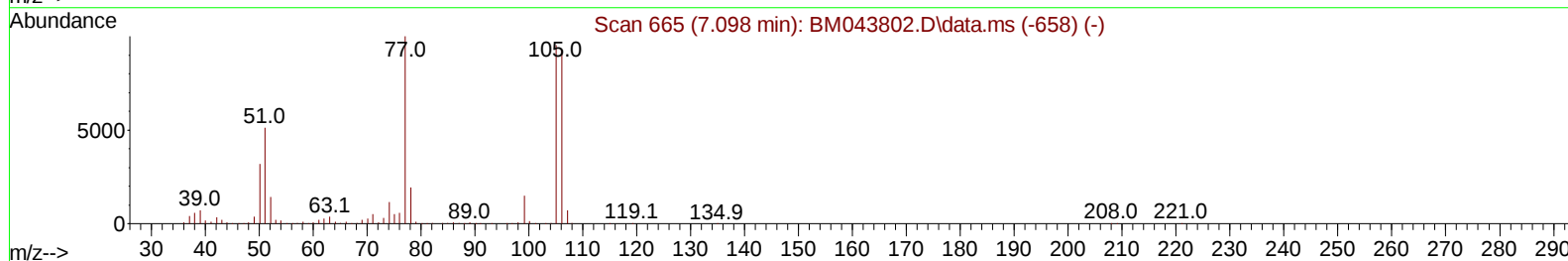
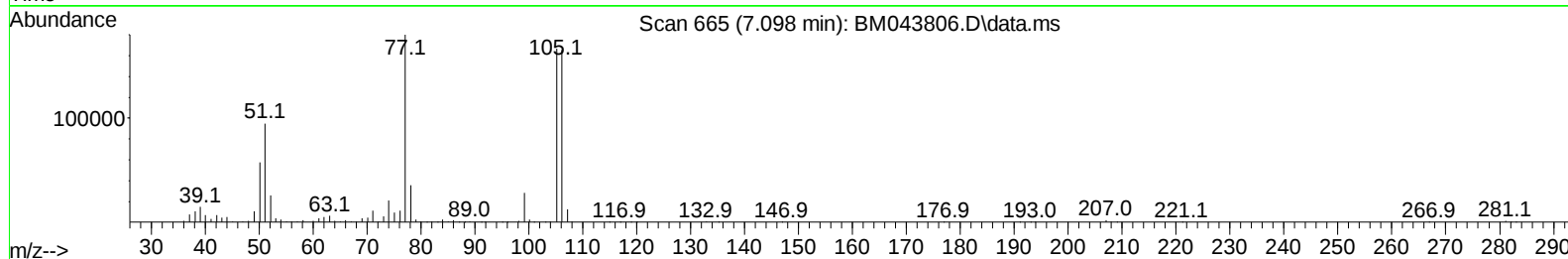
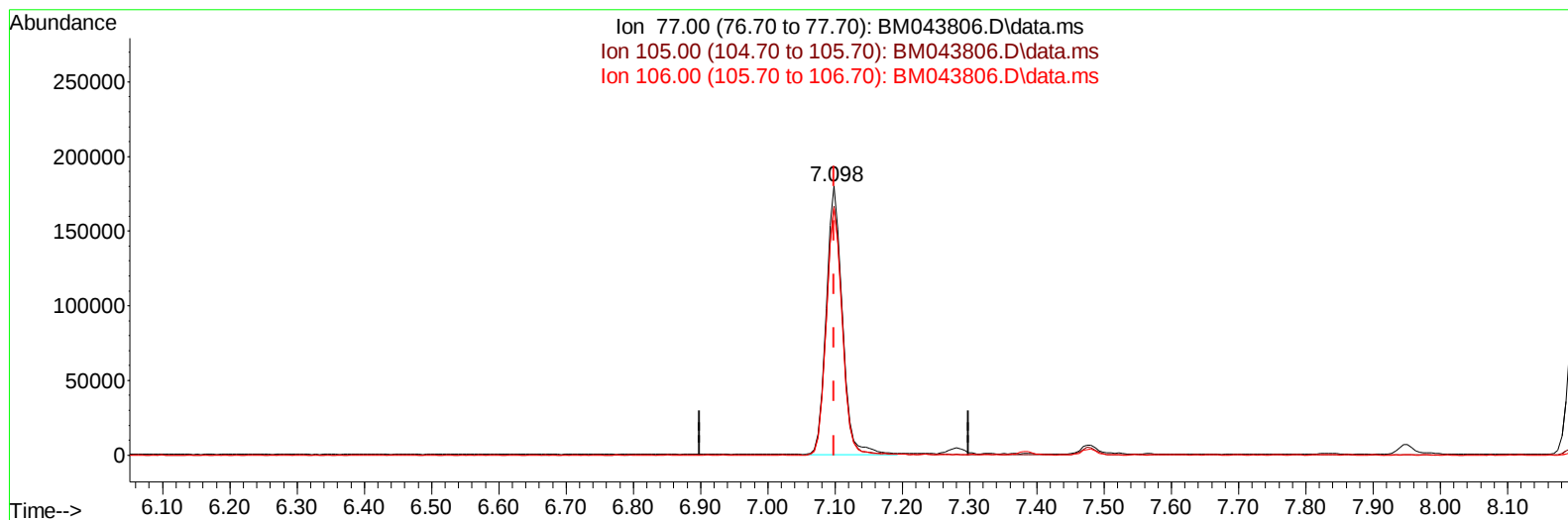
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010924\
Data File : BM043806.D
Acq On : 09 Jan 2024 17:14
Operator : MA/JU
Sample : SSTDICV020
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0B37

Manual IntegrationsAPPROVED

Quant Time: Jan 10 01:07:32 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM010924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jan 10 01:07:47 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/15/2024
Supervised By :mohammad ahmed 01/16/2024



TIC: BM043806.D\data.ms

(6) Benzaldehyde

7.098min (-0.000) 26.95 ng/u1

response 298799

Ion	Exp%	Act%
77.00	100.00	100.00
105.00	95.00	92.47
106.00	92.80	91.83
0.00	0.00	0.00

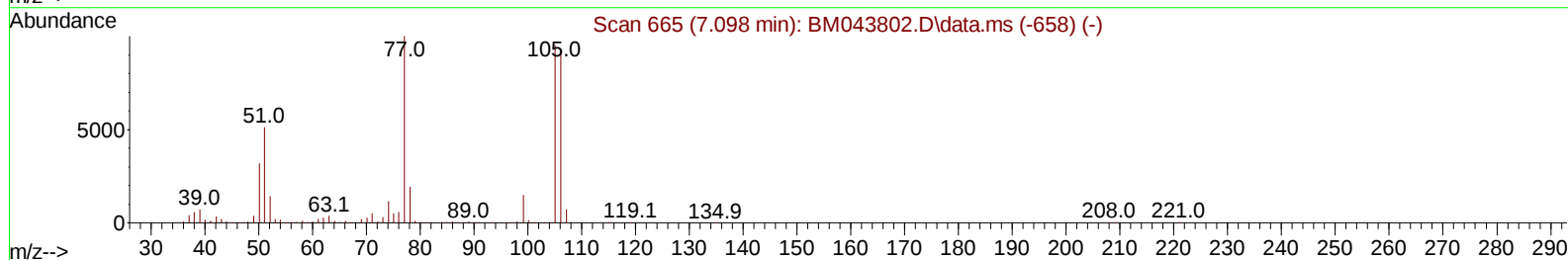
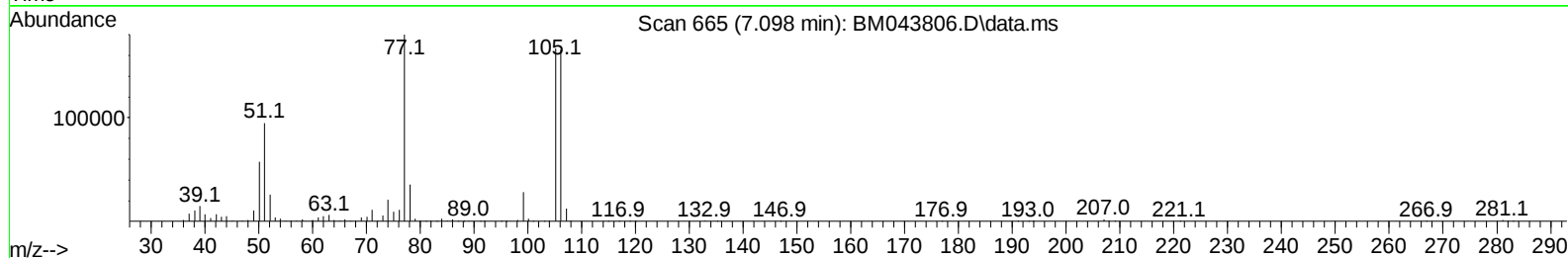
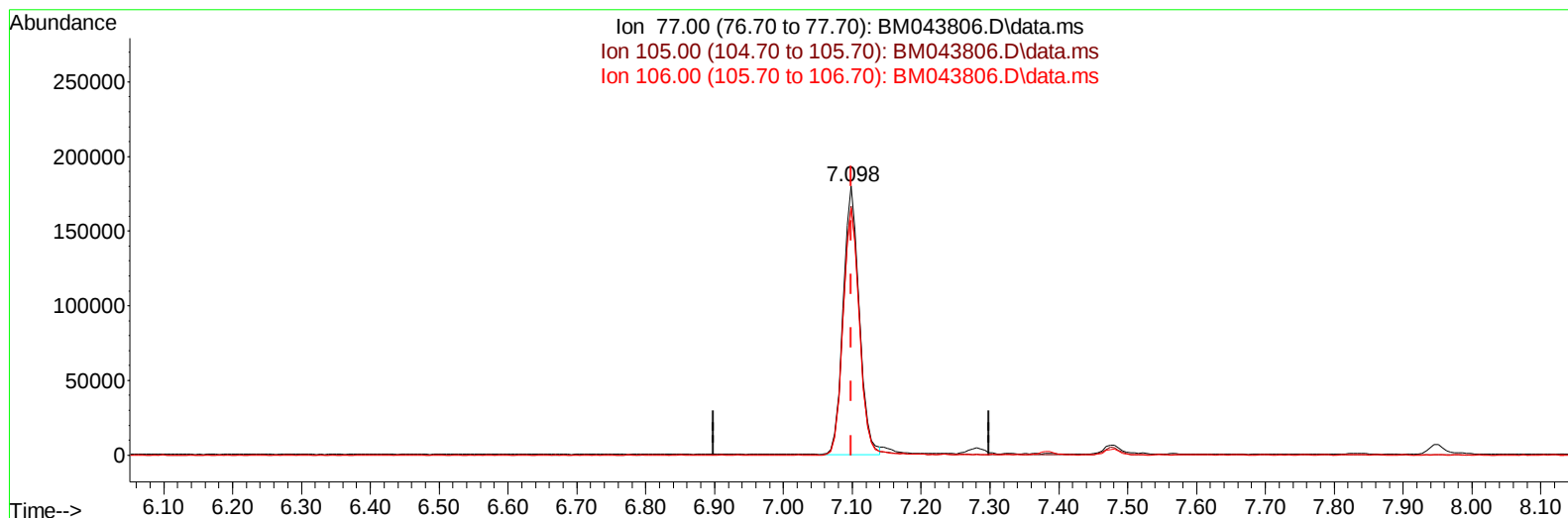
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010924\
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BM043806.D\data.ms

(6) Benzaldehyde

7.098min (-0.000) 26.41 ng/ul m

response 292848

Ion	Exp%	Act%
77.00	100.00	100.00
105.00	95.00	92.47
106.00	92.80	91.83
0.00	0.00	0.00

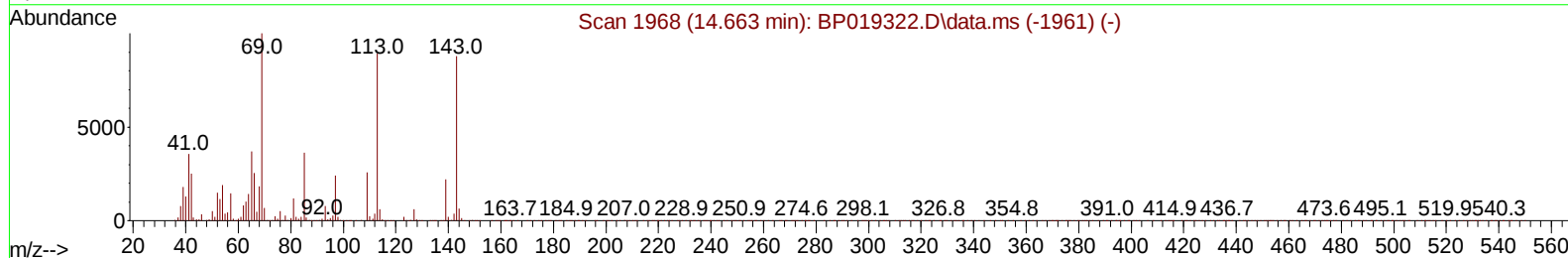
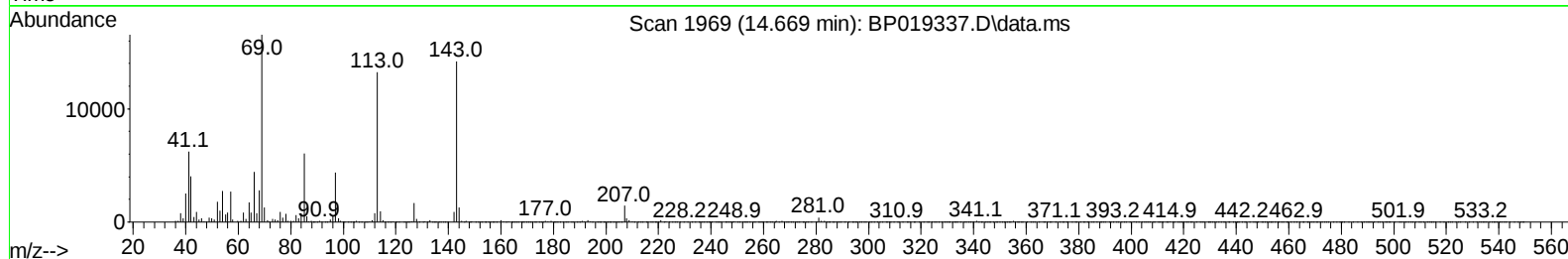
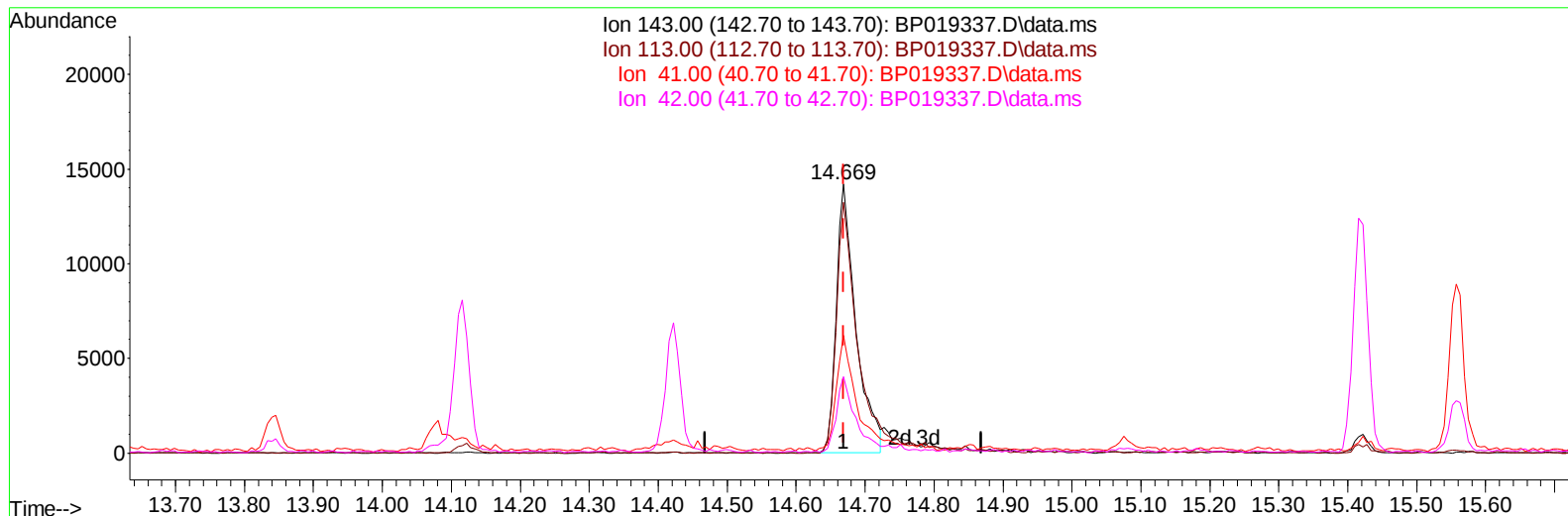
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
Data File : BP019337.D
Acq On : 12 Jan 2024 19:14
Operator : MA/JU
Sample : P1091-10
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0B37

Manual IntegrationsAPPROVED

Quant Time: Jan 12 21:38:19 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jan 09 03:05:24 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/15/2024
Supervised By :mohammad ahmed 01/16/2024



TIC: BP019337.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.000) 6.49 ng/u1

response 27995

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	93.32
41.00	39.60	43.82
42.00	25.70	28.45

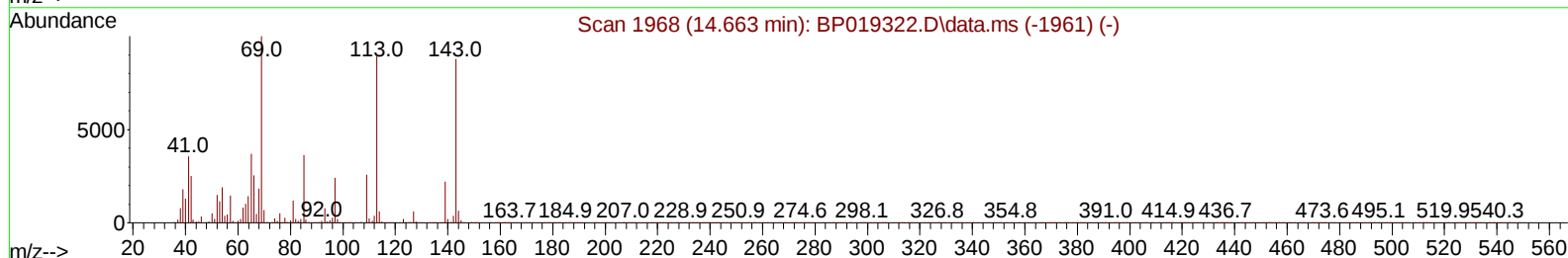
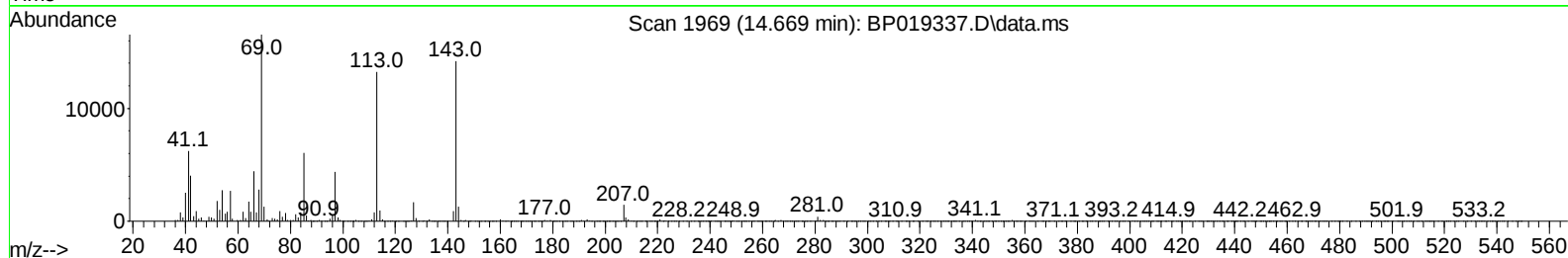
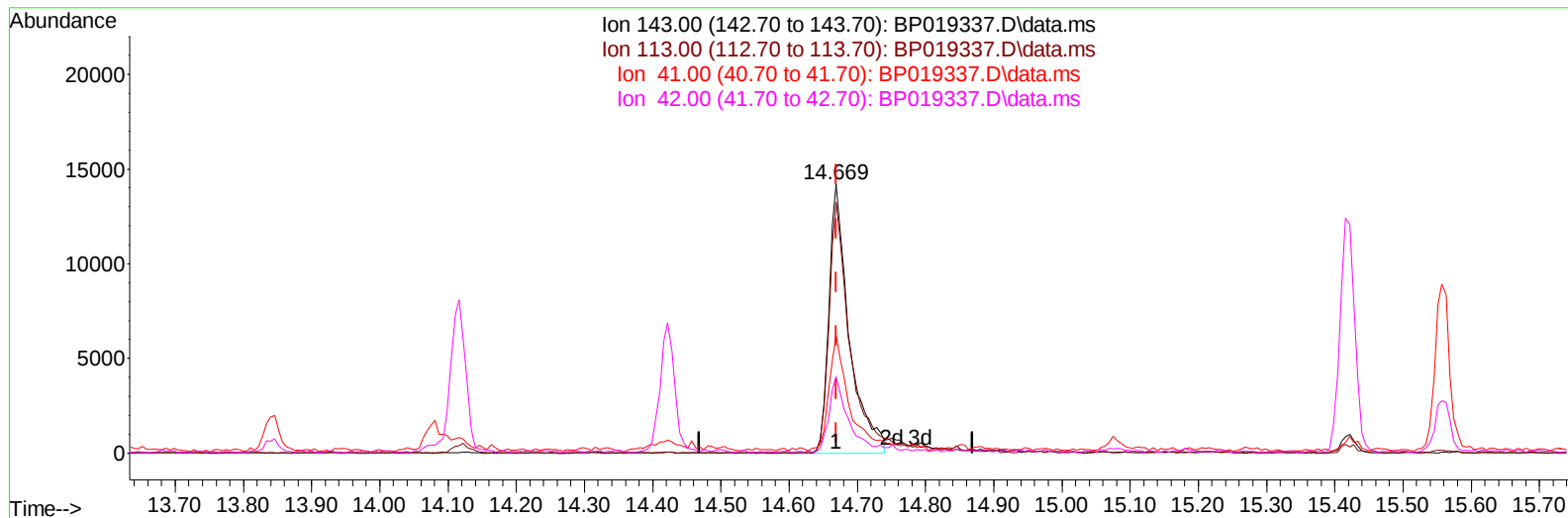
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
Data File : BP019337.D
Acq On : 12 Jan 2024 19:14
Operator : MA/JU
Sample : P1091-10
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0B37

Manual IntegrationsAPPROVED

Quant Time: Jan 12 21:38:19 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
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TIC: BP019337.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.000) 6.77 ng/ul m

response 29193

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	93.32
41.00	39.60	43.82
42.00	25.70	28.45

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
 Data File : BP019337.D
 Acq On : 12 Jan 2024 19:14
 Operator : MA/JU
 Sample : P1091-10
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B37

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/15/2024
 Supervised By :mohammad ahmed 01/16/2024

Quant Time: Jan 12 21:42:03 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 09 03:05:24 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	111523	20.000	ng/ul	0.00
20) Naphthalene-d8	10.575	136	452906	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.422	164	317884	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.181	188	808164	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240	833368	20.000	ng/ul	0.00
88) Perylene-d12	23.639	264	873377	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	11959	4.055	ng/uL	0.00
4) Pyridine-d5	3.646	84	52534	6.160	ng/ul	0.00
7) Phenol-d5	6.969	99	69767	6.545	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	202022	31.687	ng/ul	0.00
11) 2-Chlorophenol-d4	7.316	132	175590	23.981	ng/ul	0.00
15) 4-Methylphenol-d8	8.505	113	127358	15.061	ng/ul	-0.01
21) Nitrobenzene-d5	8.946	128	120199	31.546	ng/ul	0.00
24) 2-Nitrophenol-d4	9.663	143	137392	31.033	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.205	165	246716	27.584	ng/ul	0.00
31) 4-Chloroaniline-d4	10.728	131	260457	23.053	ng/ul	0.00
46) Dimethylphthalate-d6	13.846	166	926897	33.004	ng/ul	0.00
49) Acenaphthylene-d8	14.116	160	954861	31.718	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	29193m	6.769	ng/ul	0.00
60) Fluorene-d10	15.422	176	814387	35.419	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	200078	33.533	ng/ul	0.00
73) Anthracene-d10	17.281	188	1532435	38.019	ng/ul	0.00
81) Pyrene-d10	19.528	212	2009820	36.985	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.486	264	1874220	39.466	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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