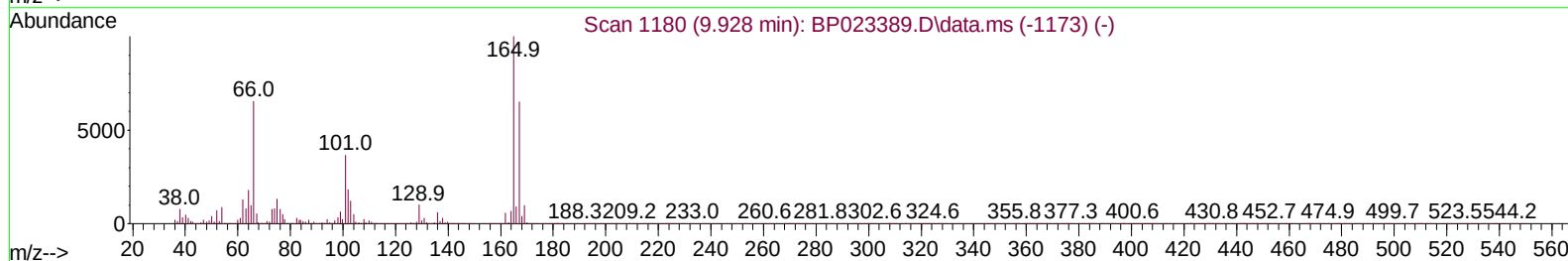
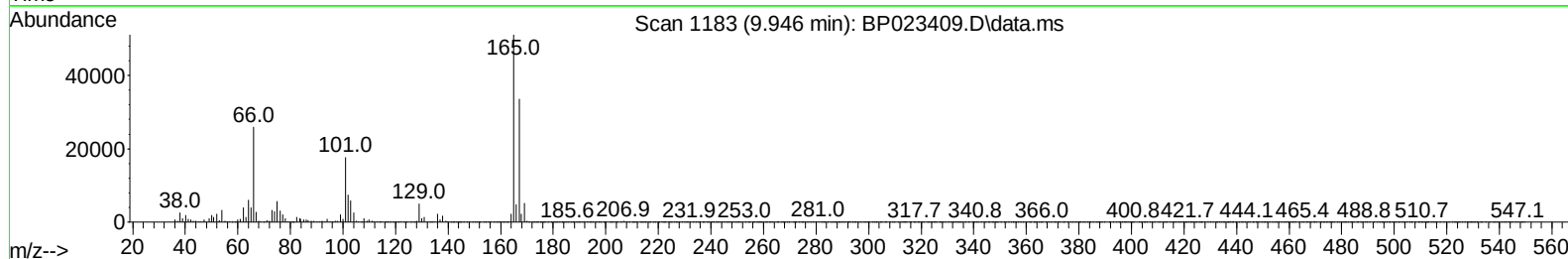
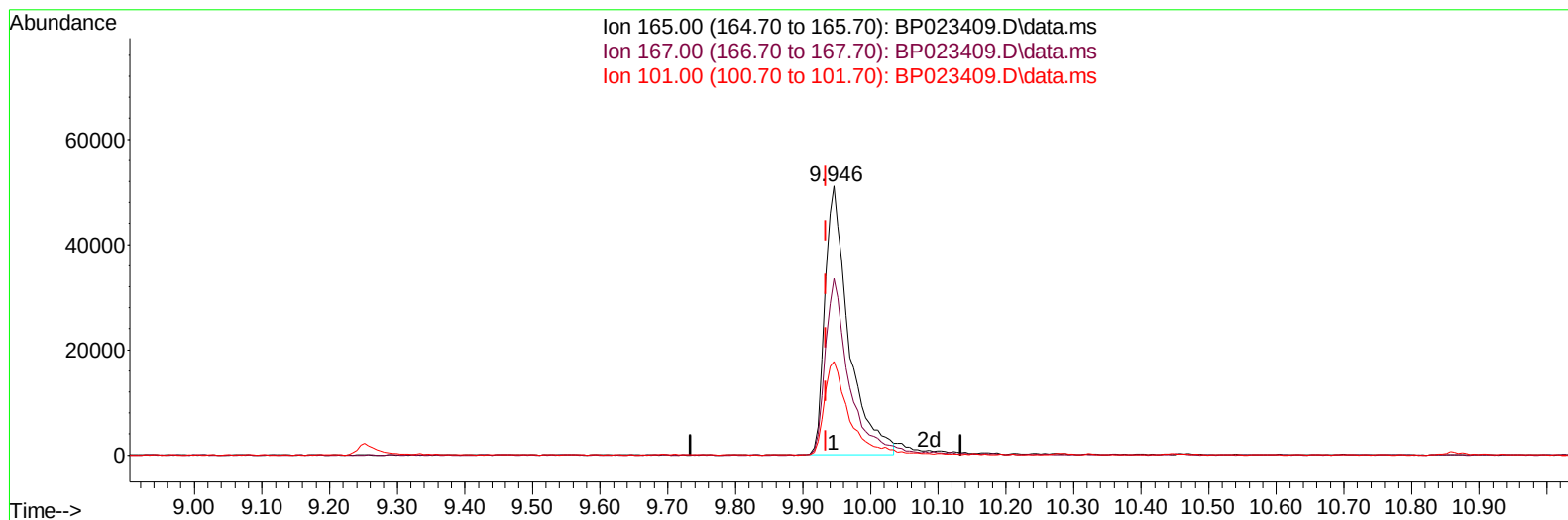


Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121224\
Data File : BP023409.D
Acq On : 12 Dec 2024 23:55
Operator : RC/JU
Sample : P5232-16
Misc :
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 13 03:44:40 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 00:30:30 2024
Response via : Initial Calibration



TIC: BP023409.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

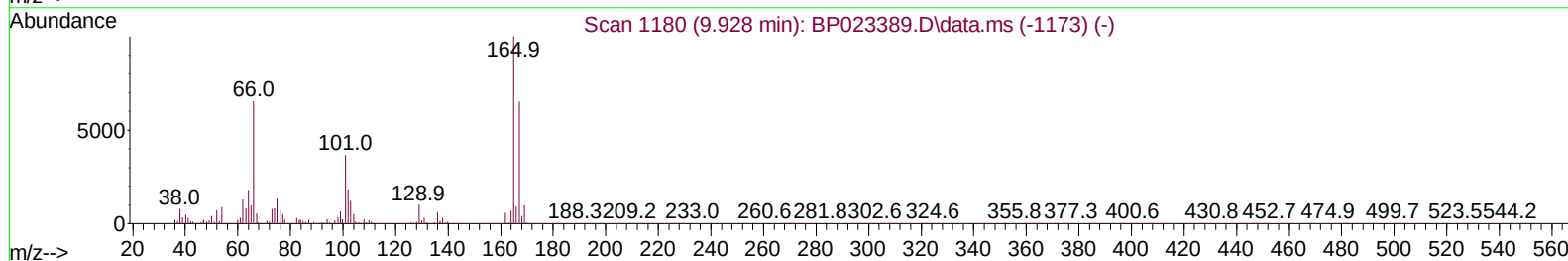
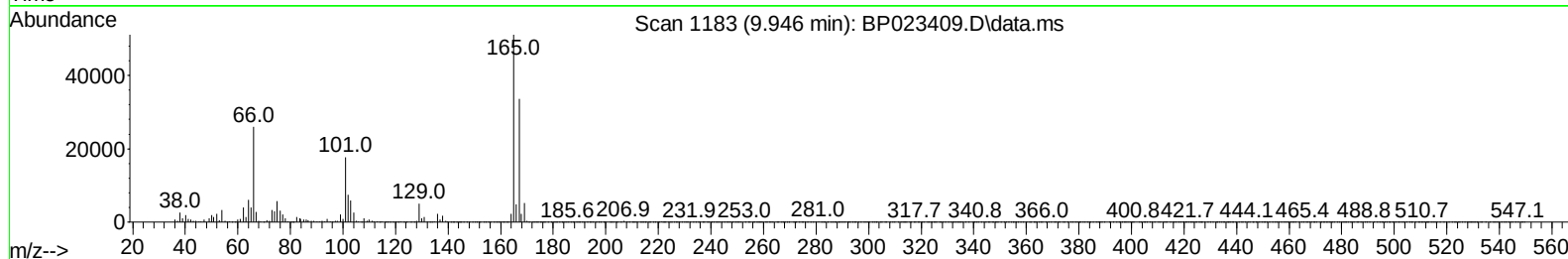
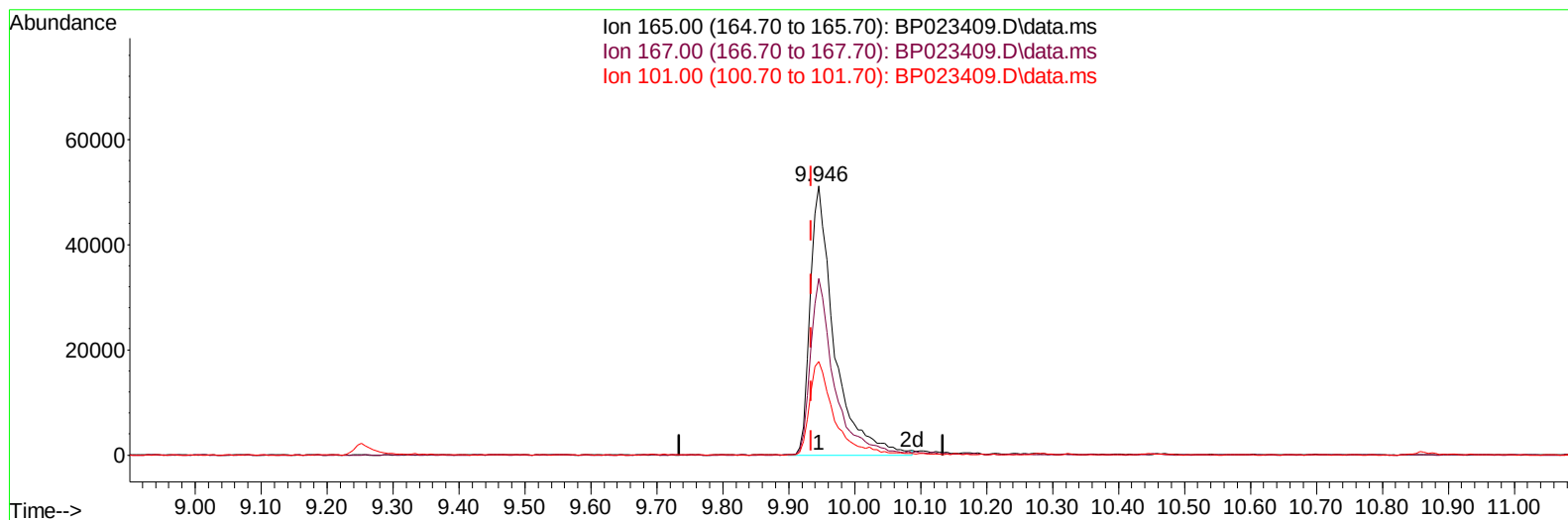
9.946min (+ 0.012) 15.15 ng/ul

response 124758

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	62.00	65.75
101.00	34.30	34.73
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121224\
Data File : BP023409.D
Acq On : 12 Dec 2024 23:55
Operator : RC/JU
Sample : P5232-16
Misc :
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 13 03:44:40 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 00:30:30 2024
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TIC: BP023409.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

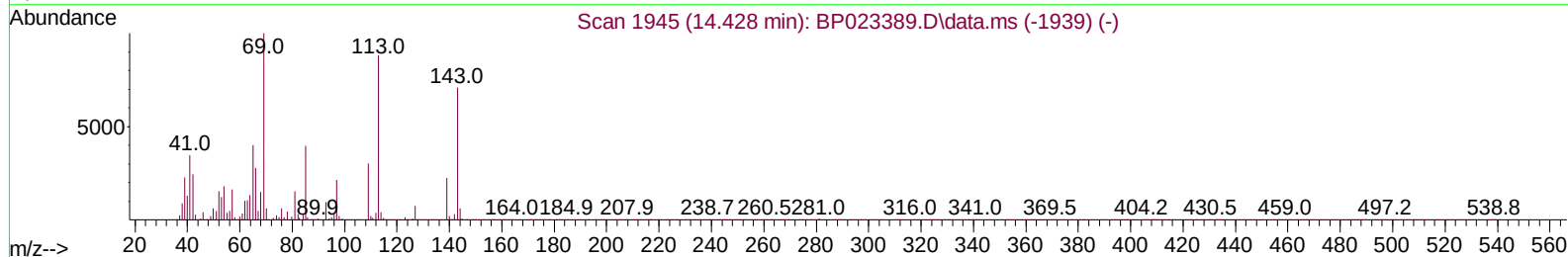
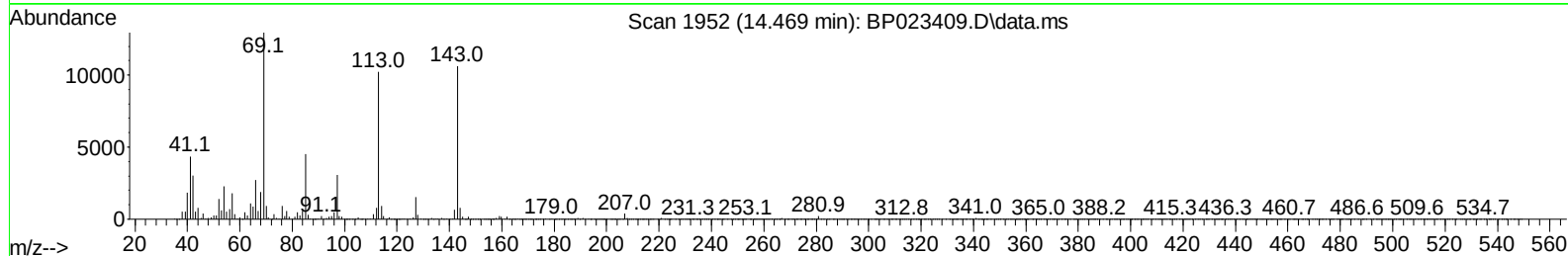
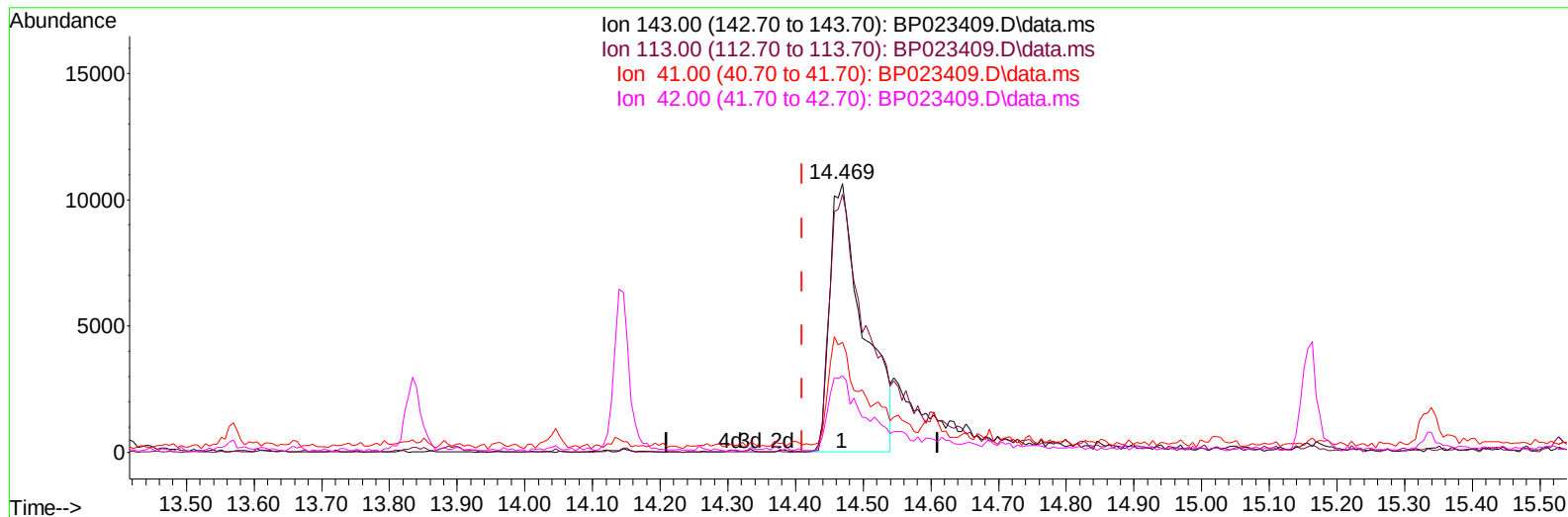
9.946min (+ 0.012) 15.77 ng/ul m

response 129868

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	62.00	65.75
101.00	34.30	34.73
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121224\
Data File : BP023409.D
Acq On : 12 Dec 2024 23:55
Operator : RC/JU
Sample : P5232-16
Misc :
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 13 03:44:40 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 00:30:30 2024
Response via : Initial Calibration



TIC: BP023409.D\data.ms

(54) 4-Nitrophenol-d4 (S)

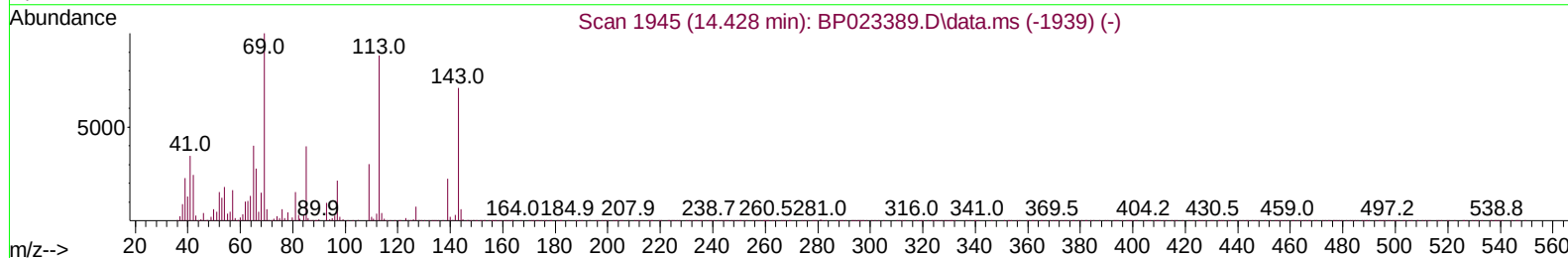
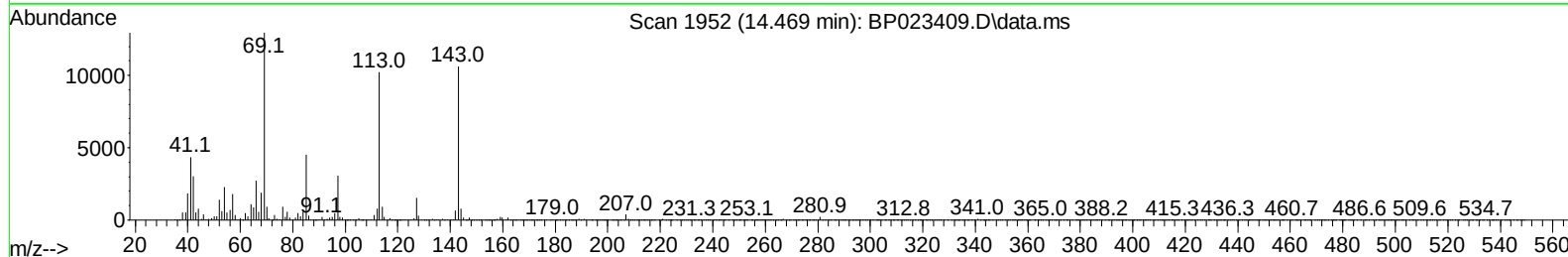
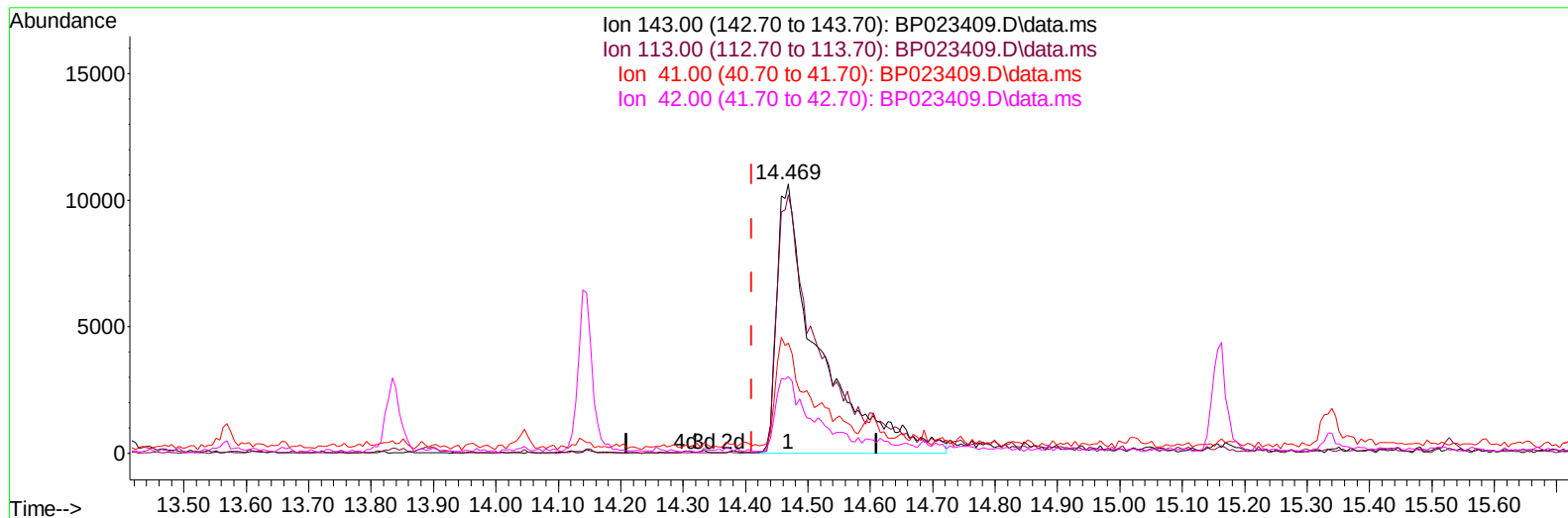
14.469min (+ 0.059) 11.50 ng/ul

response 36522

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.10	96.09
41.00	45.60	40.96
42.00	31.70	28.42

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121224\
Data File : BP023409.D
Acq On : 12 Dec 2024 23:55
Operator : RC/JU
Sample : P5232-16
Misc :
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 13 03:44:40 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 00:30:30 2024
Response via : Initial Calibration



TIC: BP023409.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.469min (+ 0.059) 15.56 ng/ul m

response 49421

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.10	96.09
41.00	45.60	40.96
42.00	31.70	28.42

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121224\
 Data File : BP023409.D
 Acq On : 12 Dec 2024 23:55
 Operator : RC/JU
 Sample : P5232-16
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 13 03:47:23 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 11 00:30:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.534	152	122210	20.000	ng/ul	0.00
20) Naphthalene-d8	10.281	136	453321	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.146	164	333940	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.951	188	764326	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	803069	20.000	ng/ul	0.00
88) Perylene-d12	24.545	264	920743	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.099	96	7922	2.387	ng/uL	0.00
4) Pyridine-d5	3.517	84	105223	12.327	ng/ul	0.00
7) Phenol-d5	6.752	99	140776	14.681	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.887	67	90469	14.954	ng/ul	0.00
11) 2-Chlorophenol-d4	7.087	132	110267	15.581	ng/ul	0.00
15) 4-Methylphenol-d8	8.263	113	63099	8.153	ng/ul	0.00
21) Nitrobenzene-d5	8.681	128	51692	15.270	ng/ul	0.00
24) 2-Nitrophenol-d4	9.393	143	61860	15.856	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.946	165	129868m	15.772	ng/ul	0.01
31) 4-Chloroaniline-d4	10.446	131	114428	12.405	ng/ul	0.00
46) Dimethylphthalate-d6	13.569	166	427726	16.499	ng/ul	0.01
49) Acenaphthylene-d8	13.834	160	376840	13.781	ng/ul	0.00
54) 4-Nitrophenol-d4	14.469	143	49421m	15.557	ng/ul	0.06
60) Fluorene-d10	15.157	176	284694	12.485	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.340	200	40269	8.718	ng/ul	0.03
73) Anthracene-d10	17.057	188	496229	14.226	ng/ul	0.00
81) Pyrene-d10	19.439	212	523841	12.485	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.339	264	528860	10.730	ng/ul	0.00

Target Compounds	Qvalue

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

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