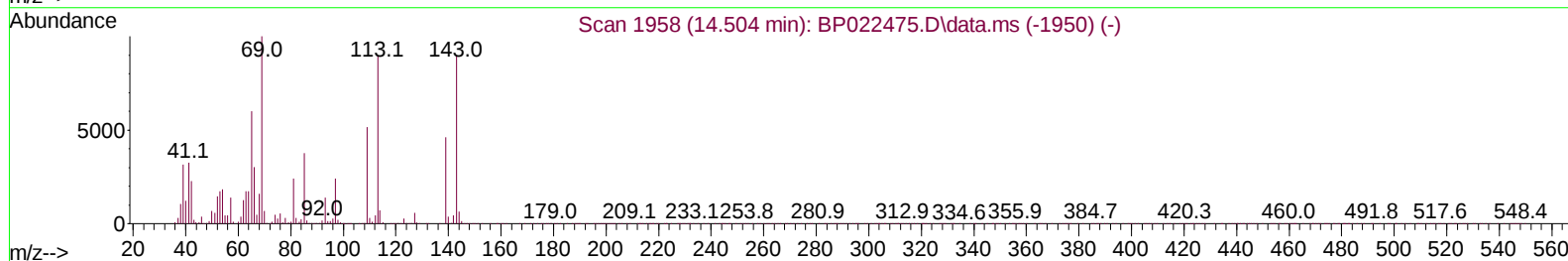
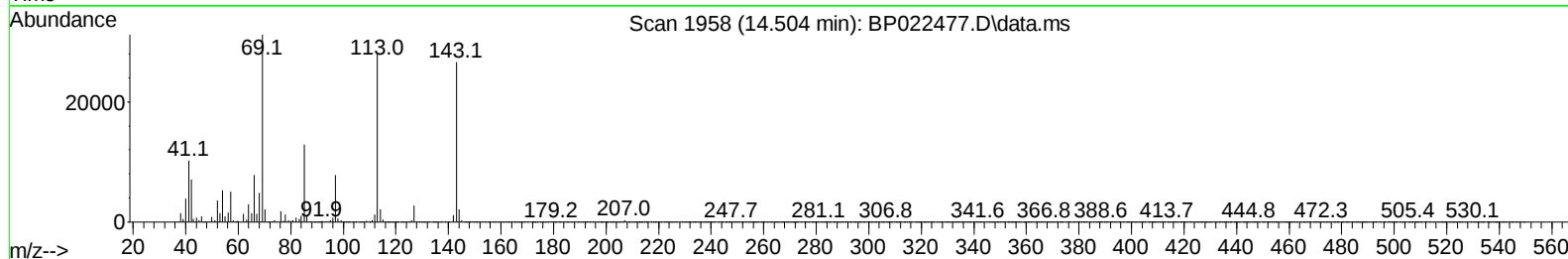
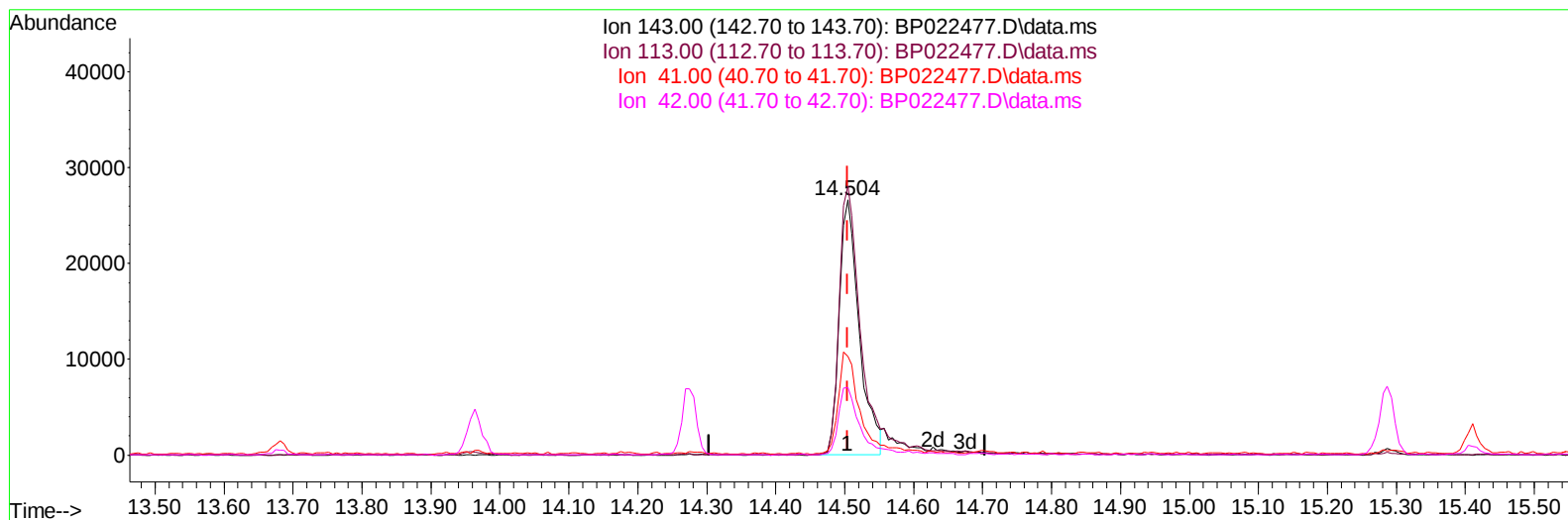


Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101924\
Data File : BP022477.D
Acq On : 19 Oct 2024 15:37
Operator : RC/JU
Sample : P4346-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Oct 21 01:37:08 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 18 11:08:24 2024
Response via : Initial Calibration



TIC: BP022477.D\data.ms

(54) 4-Nitrophenol-d4 (S)

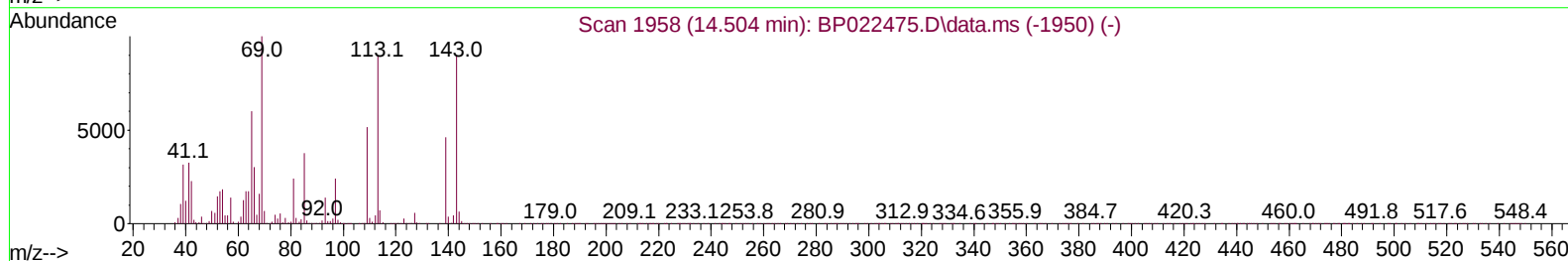
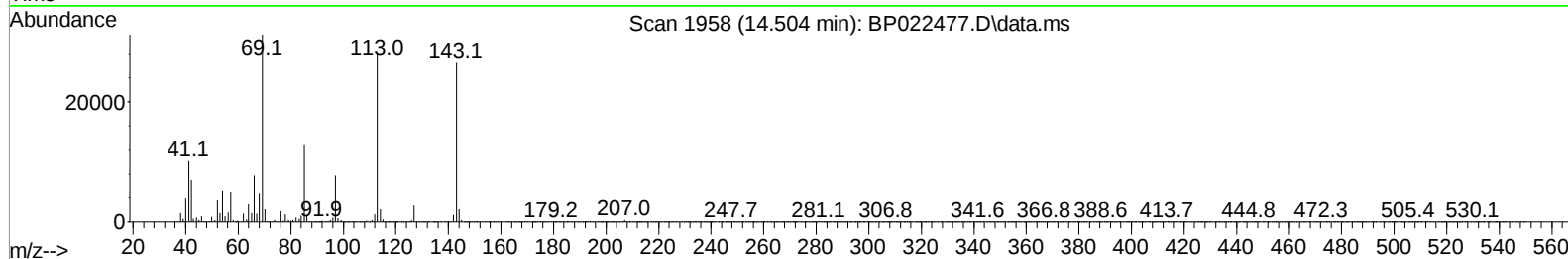
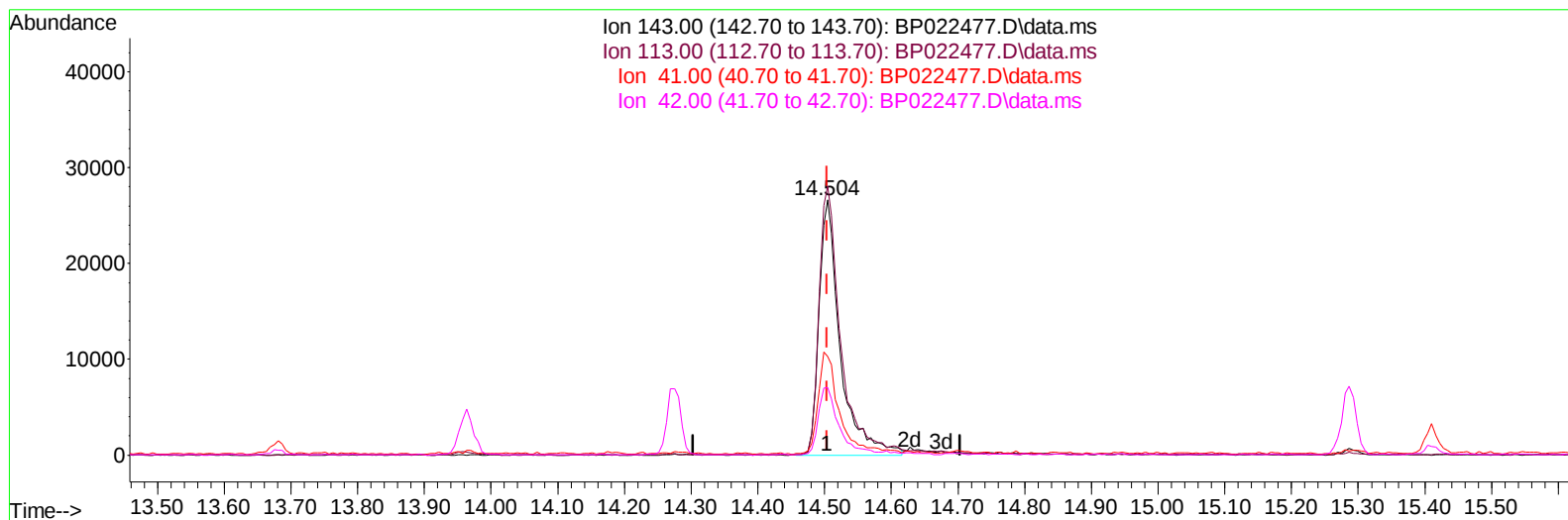
14.504min (+ 0.001) 11.93 ng/u1

response 53557

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	105.47
41.00	35.90	38.61
42.00	23.70	26.43

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101924\
Data File : BP022477.D
Acq On : 19 Oct 2024 15:37
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Sample : P4346-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Oct 21 01:37:08 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 18 11:08:24 2024
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TIC: BP022477.D\data.ms

(54) 4-Nitrophenol-d4 (S)

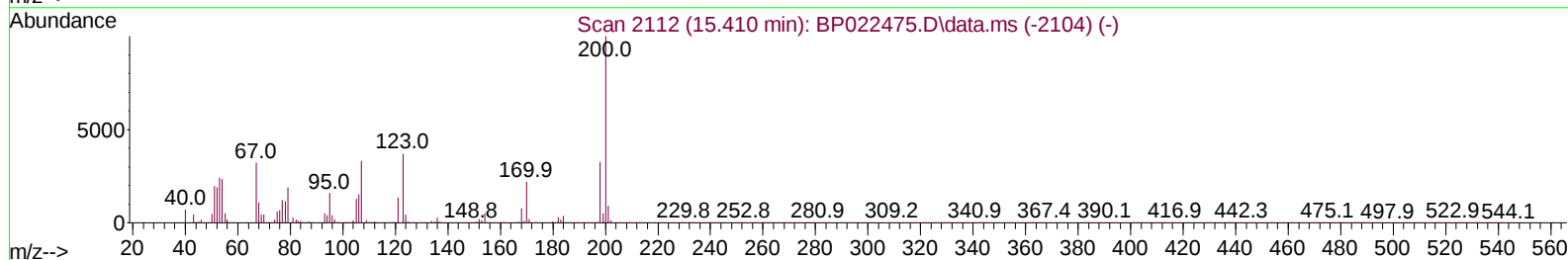
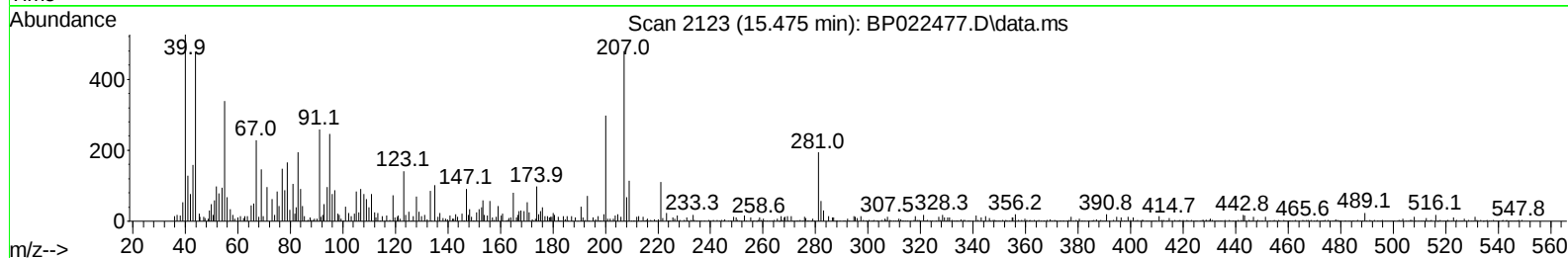
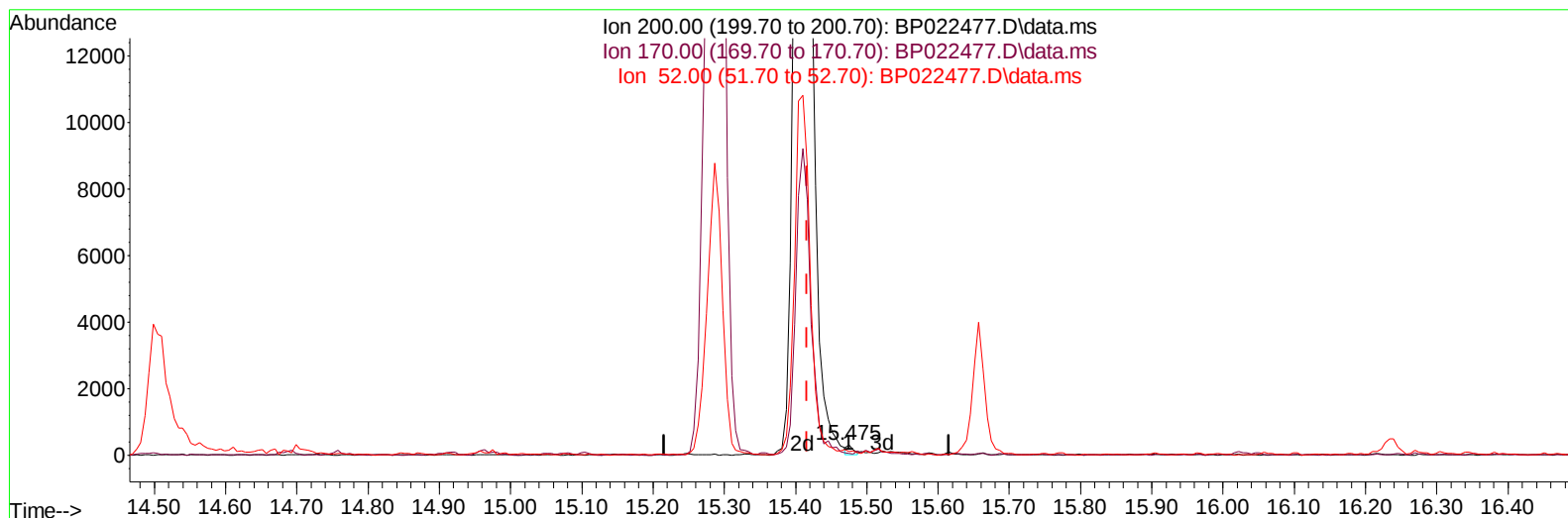
14.504min (+ 0.001) 13.02 ng/ul m

response 58464

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	105.47
41.00	35.90	38.61
42.00	23.70	26.43

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101924\
Data File : BP022477.D
Acq On : 19 Oct 2024 15:37
Operator : RC/JU
Sample : P4346-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Oct 21 01:37:08 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 18 11:08:24 2024
Response via : Initial Calibration



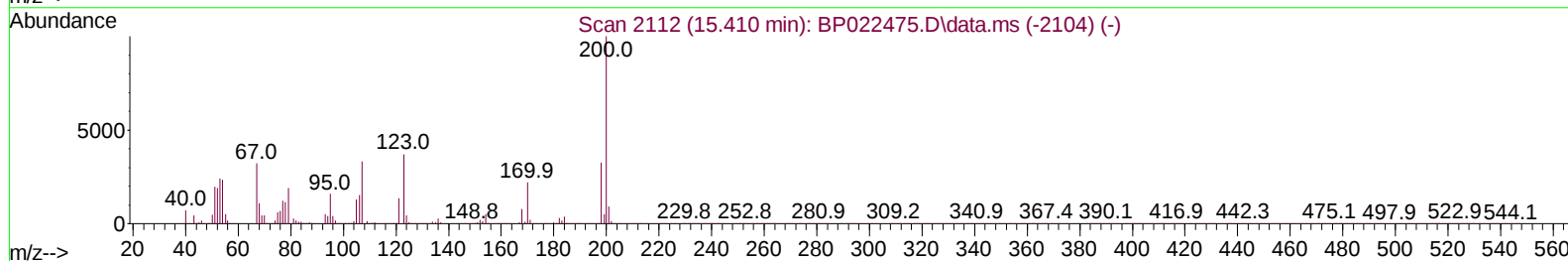
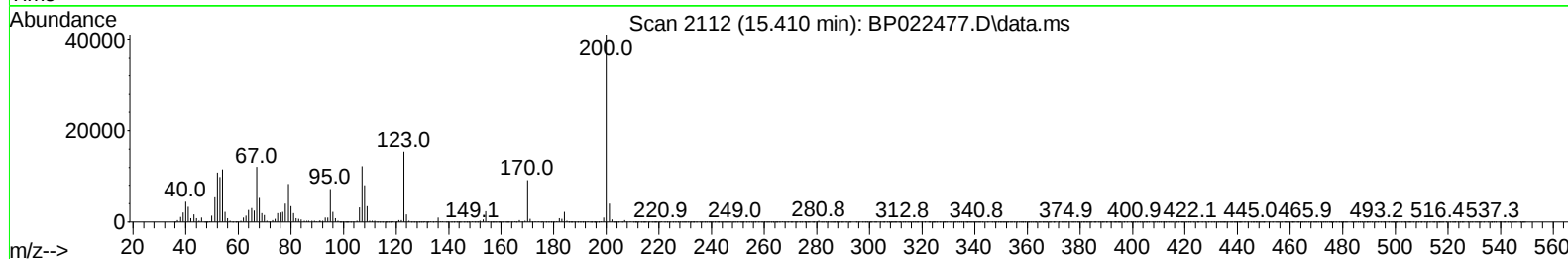
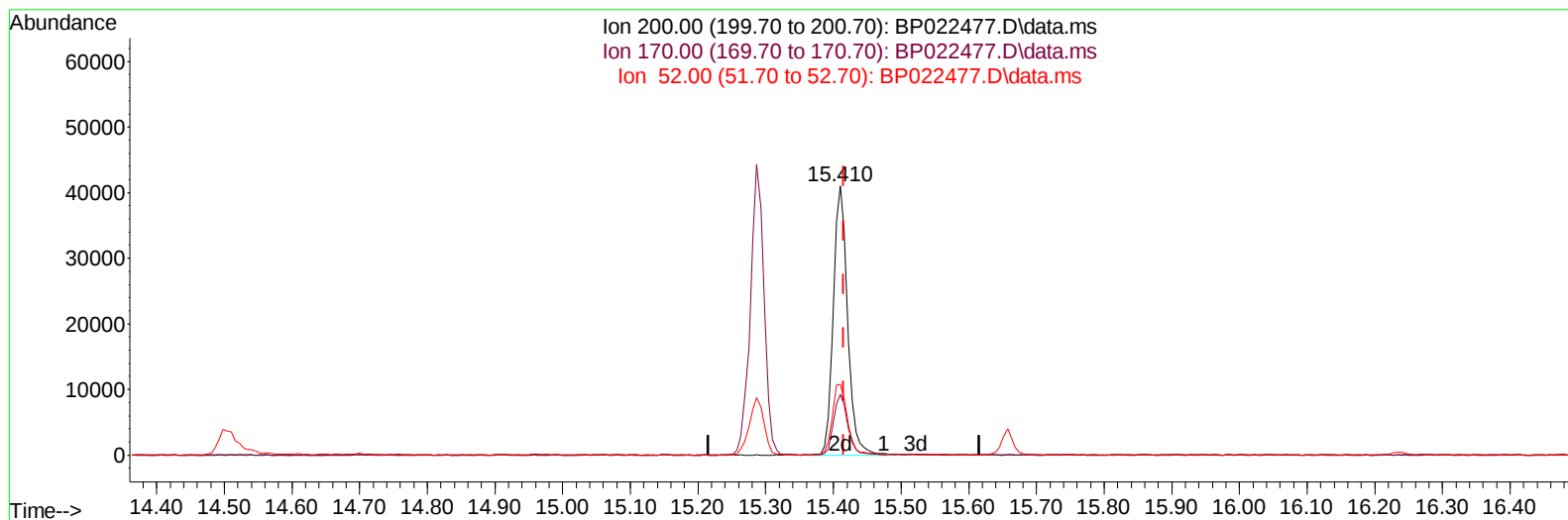
TIC: BP022477.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)**15.475min (+ 0.059) 0.03 ng/u1****response 172**

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.00	17.79
52.00	36.60	33.22
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101924\
Data File : BP022477.D
Acq On : 19 Oct 2024 15:37
Operator : RC/JU
Sample : P4346-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Oct 21 01:37:08 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 18 11:08:24 2024
Response via : Initial Calibration



TIC: BP022477.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.410min (-0.005) 11.16 ng/ul m

response 59707

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.00	22.49
52.00	36.60	26.41#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101924\
 Data File : BP022477.D
 Acq On : 19 Oct 2024 15:37
 Operator : RC/JU
 Sample : P4346-18
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Oct 21 01:41:40 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 18 11:08:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.664	152	96440	20.000	ng/ul	0.00
20) Naphthalene-d8	10.416	136	396610	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.275	164	336775	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.069	188	813653	20.000	ng/ul	-0.01
79) Chrysene-d12	21.492	240	795555	20.000	ng/ul	# 0.00
88) Perylene-d12	24.745	264	851549	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	6877	2.771	ng/uL	0.00
4) Pyridine-d5	3.611	84	100951	14.818	ng/ul	0.00
7) Phenol-d5	6.852	99	140660	17.327	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	86326	18.093	ng/ul	0.00
11) 2-Chlorophenol-d4	7.205	132	105072	18.241	ng/ul	0.00
15) 4-Methylphenol-d8	8.369	113	127651	19.703	ng/ul	0.00
21) Nitrobenzene-d5	8.805	128	55577	18.225	ng/ul	0.00
24) 2-Nitrophenol-d4	9.516	143	70602	19.997	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.058	165	143472	19.120	ng/ul	0.00
31) 4-Chloroaniline-d4	10.558	131	173818	19.603	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	565796	21.145	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	571444	20.107	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	58464m	13.024	ng/ul	0.00
60) Fluorene-d10	15.287	176	497565	21.439	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	59707m	11.158	ng/ul	0.00
73) Anthracene-d10	17.169	188	828485	21.645	ng/ul	0.00
81) Pyrene-d10	19.545	212	1031805	24.526	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.522	264	1005358	22.107	ng/ul	0.00
Target Compounds						
86) Bis(2-ethylhexyl)phtha...	21.410	149	38074	1.470	ng/ul	Qvalue 98

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

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