

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030124\
 Data File : BM044476.D
 Acq On : 01 Mar 2024 18:29
 Operator : MA/JU
 Sample : P1580-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGRF9

Quant Time: Mar 01 22:06:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 01 16:56:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.081	152	238098	20.000	ng/u1	0.03
20) Naphthalene-d8	10.916	136	973556	20.000	ng/u1	0.04
38) Acenaphthene-d10	14.745	164	562736	20.000	ng/u1	0.03
64) Phenanthrene-d10	17.498	188	1127650	20.000	ng/u1	0.02
79) Chrysene-d12	21.692	240	940451	20.000	ng/u1	0.02
88) Perylene-d12	24.233	264	1075523	20.000	ng/u1	0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.411	96	30169	3.998	ng/uL	0.01
4) Pyridine-d5	3.840	84	404193	19.612	ng/u1	0.01
7) Phenol-d5	7.240	99	500316	21.935	ng/u1	0.03
9) Bis-(2-Chloroethyl)eth...	7.416	67	333541	22.582	ng/u1	0.03
11) 2-Chlorophenol-d4	7.598	132	401969	23.399	ng/u1	0.03
15) 4-Methylphenol-d8	8.804	113	378267	21.918	ng/u1	0.03
21) Nitrobenzene-d5	9.269	128	194972	24.631	ng/u1	0.04
24) 2-Nitrophenol-d4	9.992	143	203473	25.195	ng/u1	0.03
28) 2,4-Dichlorophenol-d3	10.534	165	329828	23.152	ng/u1	0.04
31) 4-Chloroaniline-d4	11.069	131	535480	22.380	ng/u1	0.04
46) Dimethylphthalate-d6	14.169	166	1029569	23.722	ng/u1	0.02
49) Acenaphthylene-d8	14.439	160	1232109	24.675	ng/u1	0.03
54) 4-Nitrophenol-d4	14.957	143	140492	16.268	ng/u1	0.03
60) Fluorene-d10	15.739	176	909806	24.767	ng/u1	0.03
65) 4,6-Dinitro-2-methylph...	15.863	200	117479	16.790	ng/u1	0.02
73) Anthracene-d10	17.598	188	1396671	25.976	ng/u1	0.02
81) Pyrene-d10	19.886	212	1590971	28.120	ng/u1	0.02
92) Benzo(a)pyrene-d12	24.068	264	1462549	26.538	ng/u1	0.04

Target Compounds Qvalue

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

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