

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040624\
 Data File : BM044991.D
 Acq On : 07 Apr 2024 04:02
 Operator : MA/JU
 Sample : P2013-07
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Apr 08 05:06:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 08 05:01:54 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	80795	20.000	ng/ul	0.00
20) Naphthalene-d8	10.404	136	352874	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.280	164	224651	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.039	188	486160	20.000	ng/ul	0.00
79) Chrysene-d12	21.245	240	474427	20.000	ng/ul	0.00
88) Perylene-d12	23.468	264	562578	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	10891	5.034	ng/uL	0.00
4) Pyridine-d5	3.616	84	69133	11.750	ng/ul	0.00
7) Phenol-d5	6.816	99	50830	7.421	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	127932	31.350	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	147200	27.431	ng/ul	0.00
15) 4-Methylphenol-d8	8.345	113	93850	17.505	ng/ul	0.00
21) Nitrobenzene-d5	8.792	128	87964	32.453	ng/ul	0.00
24) 2-Nitrophenol-d4	9.504	143	96798	33.834	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.039	165	176047	33.562	ng/ul	0.00
31) 4-Chloroaniline-d4	10.569	131	130221	15.593	ng/ul	0.00
46) Dimethylphthalate-d6	13.704	166	565501	36.153	ng/ul	0.00
49) Acenaphthylene-d8	13.974	160	656756	35.541	ng/ul	0.00
54) 4-Nitrophenol-d4	14.533	143	15488	4.744	ng/ul	0.02
60) Fluorene-d10	15.280	176	514834	36.919	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.421	200	76955	26.697	ng/ul	0.00
73) Anthracene-d10	17.139	188	825798	37.759	ng/ul	0.00
81) Pyrene-d10	19.445	212	993348	42.522	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.327	264	1046490	38.132	ng/ul	0.00

Target Compounds Qvalue

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

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