

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101121\
 Data File : BM032451.D
 Acq On : 12 Oct 2021 10:28
 Operator : CG/JU
 Sample : M4098-16
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 YAT57

Quant Time: Oct 12 11:25:36 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM100521.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 12 09:54:27 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.607	152	264691	20.000	ng/ul	0.00
20) Naphthalene-d8	10.401	136	1210314	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.260	164	834657	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.983	188	1756223	20.000	ng/ul	0.00
79) Chrysene-d12	21.153	240	1729964	20.000	ng/ul	0.00
88) Perylene-d12	23.294	264	1734768	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.955	96	32874	4.979	ng/uL	0.00
4) Pyridine-d5	3.390	84	149250	8.003	ng/ul	0.00
7) Phenol-d5	6.801	99	188855	8.391	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.937	67	394951	30.561	ng/ul	0.00
11) 2-Chlorophenol-d4	7.148	132	443495	25.807	ng/ul	0.00
15) 4-Methylphenol-d8	8.337	113	332470	18.288	ng/ul	0.00
21) Nitrobenzene-d5	8.766	128	280358	32.831	ng/ul	0.00
24) 2-Nitrophenol-d4	9.489	143	304809	33.114	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.031	165	501827	27.387	ng/ul	0.00
31) 4-Chloroaniline-d4	10.536	131	692189	25.610	ng/ul	0.00
46) Dimethylphthalate-d6	13.666	166	1923576	33.301	ng/ul	0.00
49) Acenaphthylene-d8	13.948	160	2295108	32.707	ng/ul	0.00
54) 4-Nitrophenol-d4	14.466	143	86901	7.661	ng/ul	0.00
60) Fluorene-d10	15.248	176	1701628	34.735	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.365	200	324117	32.383	ng/ul	0.00
73) Anthracene-d10	17.083	188	2914695	38.127	ng/ul	0.00
81) Pyrene-d10	19.371	212	3303714	37.960	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.165	264	3190849	36.729	ng/ul	0.00

Target Compounds Qvalue

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

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