

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101121\
 Data File : BM032455.D
 Acq On : 12 Oct 2021 12:53
 Operator : CG/JU
 Sample : M4098-14
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 YAT45

Quant Time: Oct 12 13:27:14 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	273655	20.000	ng/ul	0.00
20) Naphthalene-d8	10.401	136	1239404	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.260	164	841118	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.983	188	1785910	20.000	ng/ul	0.00
79) Chrysene-d12	21.148	240	1750566	20.000	ng/ul	0.00
88) Perylene-d12	23.294	264	1747992	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.955	96	31740	4.650	ng/uL	0.00
4) Pyridine-d5	3.396	84	56083	2.909	ng/ul	0.01
7) Phenol-d5	6.801	99	161150	6.925	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.937	67	393777	29.472	ng/ul	0.00
11) 2-Chlorophenol-d4	7.148	132	433749	24.413	ng/ul	0.00
15) 4-Methylphenol-d8	8.337	113	301195	16.025	ng/ul	0.00
21) Nitrobenzene-d5	8.760	128	280240	32.047	ng/ul	0.00
24) 2-Nitrophenol-d4	9.490	143	300799	31.912	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.031	165	503597	26.838	ng/ul	0.00
31) 4-Chloroaniline-d4	10.531	131	527528	19.060	ng/ul	0.00
46) Dimethylphthalate-d6	13.660	166	1953543	33.560	ng/ul	0.00
49) Acenaphthylene-d8	13.948	160	2276059	32.186	ng/ul	0.00
54) 4-Nitrophenol-d4	14.466	143	65709	5.748	ng/ul	0.00
60) Fluorene-d10	15.248	176	1667988	33.787	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.366	200	305762	30.041	ng/ul	0.00
73) Anthracene-d10	17.083	188	2776266	35.713	ng/ul	0.00
81) Pyrene-d10	19.371	212	3168891	35.983	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.165	264	2957488	33.786	ng/ul	0.00

Target Compounds Qvalue

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

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