

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101121\
 Data File : BM032434.D
 Acq On : 11 Oct 2021 23:39
 Operator : CG/JU
 Sample : M4029-08
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C00R3

Quant Time: Oct 12 01:03:22 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM100521.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 11 11:06:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.613	152	182534	20.000	ng/ul	0.00
20) Naphthalene-d8	10.401	136	841139	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.260	164	601554	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.989	188	1275373	20.000	ng/ul	0.00
79) Chrysene-d12	21.153	240	1299928	20.000	ng/ul	0.00
88) Perylene-d12	23.294	264	1406052	20.000	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.955	96	9375	2.059	ng/uL	0.00
4) Pyridine-d5	3.396	84	59794	4.649	ng/ul	0.01
7) Phenol-d5	6.801	99	159279	10.262	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.937	67	95660	10.734	ng/ul	0.00
11) 2-Chlorophenol-d4	7.149	132	128467	10.840	ng/ul	0.00
15) 4-Methylphenol-d8	8.337	113	118515	9.453	ng/ul	-0.01
21) Nitrobenzene-d5	8.766	128	63266	10.660	ng/ul	0.00
24) 2-Nitrophenol-d4	9.490	143	69975	10.939	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.031	165	128533	10.093	ng/ul	0.00
31) 4-Chloroaniline-d4	10.537	131	181413	9.658	ng/ul	0.00
46) Dimethylphthalate-d6	13.666	166	479697	11.523	ng/ul	0.00
49) Acenaphthylene-d8	13.948	160	551394	10.903	ng/ul	0.00
54) 4-Nitrophenol-d4	14.466	143	52084	6.371	ng/ul	-0.01
60) Fluorene-d10	15.248	176	419161	11.872	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.366	200	38083	5.239	ng/ul	0.00
73) Anthracene-d10	17.083	188	631839	11.381	ng/ul	0.00
81) Pyrene-d10	19.371	212	794148	12.144	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.159	264	805047	11.433	ng/ul	-0.01

Target Compounds Qvalue

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

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