

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101121\
 Data File : BM032423.D
 Acq On : 11 Oct 2021 16:22
 Operator : CG/JU
 Sample : M4029-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C00R4

Quant Time: Oct 11 16:58:39 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	177330	20.000	ng/ul	0.00
20) Naphthalene-d8	10.401	136	767747	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.260	164	519794	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.989	188	1106977	20.000	ng/ul	0.00
79) Chrysene-d12	21.154	240	1105496	20.000	ng/ul	0.00
88) Perylene-d12	23.301	264	1129934	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.961	96	6989	1.580	ng/uL	0.00
4) Pyridine-d5	3.402	84	29003	2.321	ng/ul	0.02
7) Phenol-d5	6.802	99	108617	7.203	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.937	67	66311	7.659	ng/ul	0.00
11) 2-Chlorophenol-d4	7.155	132	89431	7.768	ng/ul	0.00
15) 4-Methylphenol-d8	8.343	113	74907	6.150	ng/ul	0.00
21) Nitrobenzene-d5	8.766	128	42685	7.880	ng/ul	0.00
24) 2-Nitrophenol-d4	9.490	143	44465	7.615	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.037	165	83396	7.175	ng/ul	0.00
31) 4-Chloroaniline-d4	10.537	131	115650	6.746	ng/ul	0.00
46) Dimethylphthalate-d6	13.666	166	309620	8.607	ng/ul	0.00
49) Acenaphthylene-d8	13.954	160	354730	8.117	ng/ul	0.00
54) 4-Nitrophenol-d4	14.472	143	29494	4.175	ng/ul	0.00
60) Fluorene-d10	15.248	176	267188	8.758	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.372	200	19960	3.164	ng/ul	0.00
73) Anthracene-d10	17.083	188	408326	8.474	ng/ul	0.00
81) Pyrene-d10	19.371	212	495125	8.903	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.165	264	468592	8.281	ng/ul	0.00

Target Compounds	Qvalue

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

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