

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
 Data File : BM036760.D
 Acq On : 27 Sep 2022 00:15
 Operator : CG/JU
 Sample : N4748-06
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ESQS3

Quant Time: Sep 27 00:51:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 26 23:10:22 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.998	152	394098	20.000 ng/u1	0.00
20) Naphthalene-d8	10.810	136	1699462	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.627	164	1037926	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.362	188	2014068	20.000 ng/u1	0.00
79) Chrysene-d12	21.527	240	1496401	20.000 ng/u1	0.00
88) Perylene-d12	23.962	264	1259585	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.381	96	47918	4.558 ng/uL	0.00
4) Pyridine-d5	3.816	84	206349	6.730 ng/u1	0.00
7) Phenol-d5	7.157	99	285194	8.098 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	778958	36.420 ng/u1	0.00
11) 2-Chlorophenol-d4	7.528	132	772638	30.472 ng/u1	0.00
15) 4-Methylphenol-d8	8.710	113	520094	19.986 ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	502709	43.703 ng/u1	0.00
24) 2-Nitrophenol-d4	9.892	143	468617	45.482 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	828945	34.716 ng/u1	0.00
31) 4-Chloroaniline-d4	10.945	131	1046357	26.357 ng/u1	0.00
46) Dimethylphthalate-d6	14.039	166	2909572	41.564 ng/u1	0.00
49) Acenaphthylene-d8	14.322	160	3404175	40.997 ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	84683	6.237 ng/u1	0.00
60) Fluorene-d10	15.616	176	2646571	42.188 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.727	200	302363	37.945 ng/u1	0.00
73) Anthracene-d10	17.462	188	4131146	44.861 ng/u1	0.00
81) Pyrene-d10	19.745	212	4352075	59.118 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.803	264	2898369	47.099 ng/u1	0.00

Target Compounds Qvalue

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

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