

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
 Data File : BM037728.D
 Acq On : 26 Nov 2022 11:49
 Operator : CG/JU
 Sample : N5592-08
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H0AB3

Quant Time: Nov 27 21:27:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 24 01:27:15 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.110	152	209340	20.000	ng/ul	0.00
20) Naphthalene-d8	10.933	136	1016316	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.733	164	630606	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.474	188	1265428	20.000	ng/ul	0.00
79) Chrysene-d12	21.633	240	1211989	20.000	ng/ul	0.00
88) Perylene-d12	24.127	264	1183358	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.440	96	26319	5.214	ng/uL	0.00
4) Pyridine-d5	3.875	84	156557	9.528	ng/ul	0.00
7) Phenol-d5	7.257	99	179965	8.617	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.428	67	498140	35.756	ng/ul	0.00
11) 2-Chlorophenol-d4	7.634	132	425497	29.320	ng/ul	0.00
15) 4-Methylphenol-d8	8.816	113	319111	19.201	ng/ul	0.00
21) Nitrobenzene-d5	9.280	128	290124	36.414	ng/ul	0.00
24) 2-Nitrophenol-d4	10.010	143	272146	34.665	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.545	165	461661	32.134	ng/ul	0.00
31) 4-Chloroaniline-d4	11.063	131	506067	21.227	ng/ul	-0.01
46) Dimethylphthalate-d6	14.145	166	1866037	41.776	ng/ul	0.00
49) Acenaphthylene-d8	14.433	160	2049174	38.867	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	81300	8.263	ng/ul	-0.01
60) Fluorene-d10	15.721	176	1609136	41.134	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.833	200	210519	30.926	ng/ul	0.00
73) Anthracene-d10	17.574	188	2806585	47.428	ng/ul	0.00
81) Pyrene-d10	19.850	212	3208176	52.089	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.968	264	2948445	48.406	ng/ul	0.00

Target Compounds Qvalue

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

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