

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112822\
 Data File : BM037767.D
 Acq On : 28 Nov 2022 19:43
 Operator : CG/JU
 Sample : N5710-13
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Nov 28 22:29:57 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 28 22:24:47 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.104	152	219539	20.000	ng/ul	0.00
20) Naphthalene-d8	10.928	136	1020058	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.733	164	634987	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.468	188	1311791	20.000	ng/ul	0.00
79) Chrysene-d12	21.633	240	1271026	20.000	ng/ul	0.00
88) Perylene-d12	24.127	264	1098804	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	31792	6.006	ng/uL	0.00
4) Pyridine-d5	3.875	84	179473	10.416	ng/ul	0.00
7) Phenol-d5	7.257	99	667916	30.495	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.428	67	456773	31.263	ng/ul	0.00
11) 2-Chlorophenol-d4	7.634	132	504263	33.133	ng/ul	0.00
15) 4-Methylphenol-d8	8.816	113	531814	30.513	ng/ul	0.00
21) Nitrobenzene-d5	9.281	128	275346	34.432	ng/ul	0.00
24) 2-Nitrophenol-d4	10.004	143	284631	36.122	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.540	165	471168	32.675	ng/ul	0.00
31) 4-Chloroaniline-d4	11.063	131	458153	19.147	ng/ul	0.00
46) Dimethylphthalate-d6	14.139	166	1485510	33.028	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	1822974	34.338	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	278855	28.147	ng/ul	0.00
60) Fluorene-d10	15.722	176	1349405	34.256	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.827	200	176460	25.006	ng/ul	0.00
73) Anthracene-d10	17.568	188	2119853	34.557	ng/ul	0.00
81) Pyrene-d10	19.851	212	2404923	37.233	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.968	264	1944775	34.385	ng/ul	0.00

Target Compounds Qvalue

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

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