

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP113022\
 Data File : BP012882.D
 Acq On : 30 Nov 2022 16:39
 Operator : CG/JU
 Sample : N5725-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHB55

Quant Time: Nov 30 21:56:13 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 29 22:53:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.011	152	613038	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	2808622	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.645	164	1954074	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.404	188	4069898	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	2998850	20.000	ng/ul	0.00
88) Perylene-d12	23.998	264	2479514	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	56425	3.944	ng/uL	0.00
4) Pyridine-d5	3.828	84	148805	3.476	ng/ul	0.01
7) Phenol-d5	7.158	99	257988	5.121	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	847326	27.297	ng/ul	0.00
11) 2-Chlorophenol-d4	7.534	132	769196	20.039	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113	511383	12.600	ng/ul	0.00
21) Nitrobenzene-d5	9.175	128	536430	31.205	ng/ul	0.00
24) 2-Nitrophenol-d4	9.899	143	491956	26.952	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	967079	22.707	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	826032	14.164	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166	4076052	30.305	ng/ul	0.00
49) Acenaphthylene-d8	14.340	160	4417904	28.151	ng/ul	0.00
54) 4-Nitrophenol-d4	14.822	143	61127	2.549	ng/ul	0.00
60) Fluorene-d10	15.640	176	3677013	31.216	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	326271	16.773	ng/ul	0.00
73) Anthracene-d10	17.498	188	5951896	33.222	ng/ul	0.00
81) Pyrene-d10	19.728	212	6314431	36.238	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.833	264	4032157	33.107	ng/ul	0.00

Target Compounds Qvalue

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

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