

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP113022\
 Data File : BP012880.D
 Acq On : 30 Nov 2022 15:31
 Operator : CG/JU
 Sample : N5811-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGCG9

Quant Time: Nov 30 21:55:24 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	8.010	152	585155	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	2557949	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.645	164	1606369	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.398	188	2981578	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	2362550	20.000	ng/ul	0.00
88) Perylene-d12	23.998	264	2505024	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	58070	4.253	ng/uL	0.00
4) Pyridine-d5	3.822	84	210484	5.151	ng/ul	0.00
7) Phenol-d5	7.158	99	275147	5.722	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	827391	27.925	ng/ul	0.00
11) 2-Chlorophenol-d4	7.534	132	805049	21.972	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113	502212	12.963	ng/ul	0.00
21) Nitrobenzene-d5	9.175	128	491062	31.366	ng/ul	0.00
24) 2-Nitrophenol-d4	9.899	143	507580	30.533	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	923140	23.800	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	1115179	20.995	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166	3054414	27.624	ng/ul	0.00
49) Acenaphthylene-d8	14.339	160	3574889	27.710	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	46994	2.384	ng/ul	0.00
60) Fluorene-d10	15.639	176	2715941	28.048	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	305650	21.448	ng/ul	0.00
73) Anthracene-d10	17.498	188	4095962	31.208	ng/ul	0.00
81) Pyrene-d10	19.727	212	4164280	30.335	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.833	264	3744792	30.434	ng/ul	0.00

Target Compounds Qvalue

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

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