

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
 Data File : BP019168.D
 Acq On : 29 Dec 2023 16:23
 Operator : MA/JU
 Sample : 06020-06
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH7W7

Quant Time: Dec 29 18:19:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 03:48:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	202335	20.000	ng/ul	0.00
20) Naphthalene-d8	10.598	136	759307	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.445	164	466134	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	1046778	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	1137878	20.000	ng/ul	0.00
88) Perylene-d12	23.668	264	1368993	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	26328	4.444	ng/uL	0.00
4) Pyridine-d5	3.658	84	157463	10.589	ng/ul	0.00
7) Phenol-d5	6.981	99	133718	7.369	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	334536	31.869	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	354125	24.209	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	228421	16.170	ng/ul	0.00
21) Nitrobenzene-d5	8.963	128	196314	29.909	ng/ul	0.00
24) 2-Nitrophenol-d4	9.687	143	200152	28.215	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	387495	27.051	ng/ul	0.00
31) 4-Chloroaniline-d4	10.745	131	483369	27.434	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	1305905	31.718	ng/ul	0.00
49) Acenaphthylene-d8	14.139	160	1431085	32.034	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	37894	5.450	ng/ul	0.00
60) Fluorene-d10	15.439	176	1168665	32.856	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	172571	25.148	ng/ul	0.00
73) Anthracene-d10	17.298	188	1969023	36.019	ng/ul	0.00
81) Pyrene-d10	19.545	212	2505727	36.133	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.515	264	2853489	36.263	ng/ul	0.00

Target Compounds Qvalue

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

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