

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
 Data File : BP019070.D
 Acq On : 23 Dec 2023 09:36
 Operator : MA/JU
 Sample : 05869-03
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B28

Quant Time: Dec 24 23:58:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 21 04:46:12 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	189457	20.000	ng/ul	0.00
20) Naphthalene-d8	10.604	136	700505	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.451	164	459261	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	1095346	20.000	ng/ul	-0.01
79) Chrysene-d12	21.292	240	1176191	20.000	ng/ul	0.00
88) Perylene-d12	23.651	264	1362265	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	17932	3.225	ng/uL	0.00
4) Pyridine-d5	3.669	84	129431	8.691	ng/ul	0.00
7) Phenol-d5	6.993	99	324357	17.060	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	181956	16.147	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.346	132	250535	16.887	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113	254654	16.550	ng/ul	0.00
21) Nitrobenzene-d5	8.975	128	117353	18.822	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.699	143	131146	20.289	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	260798	19.731	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.751	131	322166	18.651	ng/ul	-0.01
46) Dimethylphthalate-d6	13.869	166	876024	21.372	ng/ul	0.00
49) Acenaphthylene-d8	14.139	160	939090	20.806	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.669	143	146338	21.750	ng/ul	0.00
60) Fluorene-d10	15.439	176	772178	22.443	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.569	200	121175	18.450	ng/ul	0.00
73) Anthracene-d10	17.298	188	1295565	22.499	ng/ul	0.00
81) Pyrene-d10	19.539	212	1650822	18.031	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.498	264	1773238	22.325	ng/ul	0.00

Target Compounds Qvalue

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

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