

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
 Data File : BP019063.D
 Acq On : 23 Dec 2023 05:39
 Operator : MA/JU
 Sample : 05835-12
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AQ9

Quant Time: Dec 25 00:03:57 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	157567	20.000	ng/u1	0.00
20) Naphthalene-d8	10.610	136	614604	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.451	164	404487	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.198	188	957888	20.000	ng/u1	-0.01
79) Chrysene-d12	21.292	240	1004518	20.000	ng/u1	0.00
88) Perylene-d12	23.651	264	1154263	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	26584	5.749	ng/uL	0.00
4) Pyridine-d5	3.670	84	376166	30.372	ng/u1	0.00
7) Phenol-d5	6.993	99	496565	31.403	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	280591	29.940	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.346	132	389468	31.565	ng/u1	0.00
15) 4-Methylphenol-d8	8.534	113	408133	31.894	ng/u1	0.00
21) Nitrobenzene-d5	8.981	128	192918	35.267	ng/u1	0.00
24) 2-Nitrophenol-d4	9.699	143	222497	39.232	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	425113	36.658	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.752	131	552061	36.427	ng/u1	-0.01
46) Dimethylphthalate-d6	13.869	166	1384451	38.350	ng/u1	0.00
49) Acenaphthylene-d8	14.145	160	1547576	38.931	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	243081	41.021	ng/u1	0.00
60) Fluorene-d10	15.445	176	1233873	40.718	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	218237	37.997	ng/u1	0.00
73) Anthracene-d10	17.298	188	2027901	40.270	ng/u1	0.00
81) Pyrene-d10	19.539	212	2566164	32.818	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.498	264	2747060	40.817	ng/u1	0.00

Target Compounds Qvalue

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

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