

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
 Data File : BP019067.D
 Acq On : 23 Dec 2023 07:54
 Operator : MA/JU
 Sample : 05869-08
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AT4

Quant Time: Dec 24 23:58:00 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	192735	20.000	ng/u1	0.00
20) Naphthalene-d8	10.610	136	739081	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.451	164	488693	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.198	188	1148899	20.000	ng/u1	-0.01
79) Chrysene-d12	21.292	240	1194905	20.000	ng/u1	0.00
88) Perylene-d12	23.651	264	1383658	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	30347	5.366	ng/uL	0.00
4) Pyridine-d5	3.670	84	422110	27.863	ng/u1	0.00
7) Phenol-d5	6.993	99	542492	28.048	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	303015	26.433	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.346	132	420791	27.881	ng/u1	0.00
15) 4-Methylphenol-d8	8.534	113	437600	27.957	ng/u1	0.00
21) Nitrobenzene-d5	8.975	128	207216	31.501	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.699	143	239593	35.131	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	454615	32.600	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.752	131	607969	33.359	ng/u1	-0.01
46) Dimethylphthalate-d6	13.869	166	1467603	33.648	ng/u1	0.00
49) Acenaphthylene-d8	14.140	160	1650560	34.367	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.663	143	257959	36.031	ng/u1	0.00
60) Fluorene-d10	15.445	176	1311024	35.809	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	233635	33.915	ng/u1	0.00
73) Anthracene-d10	17.298	188	2153475	35.654	ng/u1	0.00
81) Pyrene-d10	19.539	212	2694401	28.968	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.498	264	2873324	35.615	ng/u1	0.00

Target Compounds	Qvalue

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

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