

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
 Data File : BP018389.D
 Acq On : 26 Nov 2023 03:45
 Operator : MA/JU
 Sample : 05261-12
 Misc :
 ALS Vial : 73 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 27 00:07:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 24 22:02:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	473984	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	1869834	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.498	164	1062467	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	1850565	20.000	ng/ul	0.00
79) Chrysene-d12	21.357	240	1618338	20.000	ng/ul	0.00
88) Perylene-d12	23.804	264	1971593	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	51446	4.065	ng/uL	0.00
4) Pyridine-d5	3.687	84	681392	19.593	ng/ul	0.00
7) Phenol-d5	7.028	99	962623	22.223	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	552727	21.338	ng/ul	0.00
11) 2-Chlorophenol-d4	7.393	132	804906	22.131	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	781583	22.358	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	380525	25.055	ng/ul	0.00
24) 2-Nitrophenol-d4	9.746	143	371541	26.489	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	776643	22.951	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	1033637	22.648	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	2262262	23.929	ng/ul	-0.01
49) Acenaphthylene-d8	14.187	160	2588159	24.026	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	233361	17.881	ng/ul	0.00
60) Fluorene-d10	15.487	176	1826025	23.093	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	153493	19.484	ng/ul	0.00
73) Anthracene-d10	17.345	188	2442660	24.695	ng/ul	0.00
81) Pyrene-d10	19.586	212	2568215	19.940	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	2896374	25.339	ng/ul	0.02

Target Compounds Qvalue

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

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