

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
 Data File : BP018407.D
 Acq On : 27 Nov 2023 18:55
 Operator : MA/JU
 Sample : 05341-16
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 27 21:58:13 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 27 21:50:01 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	464688	20.000	ng/ul	0.00
20) Naphthalene-d8	10.657	136	1719041	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.492	164	891108	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	1709982	20.000	ng/ul	0.00
79) Chrysene-d12	21.374	240	1786269	20.000	ng/ul	0.00
88) Perylene-d12	23.892	264	2081078	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	46879	3.779	ng/uL	0.00
4) Pyridine-d5	3.687	84	237550	6.967	ng/ul	0.00
7) Phenol-d5	7.022	99	178510	4.203	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	600863	23.661	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	625748	17.549	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	346129	10.099	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	382262	27.377	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	335947	26.052	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	649939	20.891	ng/ul	0.00
31) 4-Chloroaniline-d4	10.798	131	820257	19.549	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	2119995	26.737	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160	2395374	26.513	ng/ul	0.00
54) 4-Nitrophenol-d4	14.686	143	33613	3.071	ng/ul	0.00
60) Fluorene-d10	15.486	176	1796989	27.095	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	166033	22.809	ng/ul	0.00
73) Anthracene-d10	17.345	188	2875482	31.461	ng/ul	0.00
81) Pyrene-d10	19.592	212	3505309	24.657	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.721	264	3883489	32.188	ng/ul	0.02

Target Compounds Qvalue

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

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