

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
 Data File : BP018422.D
 Acq On : 28 Nov 2023 03:57
 Operator : MA/JU
 Sample : 05332-20
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YCGC7

Quant Time: Nov 28 05:59:52 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 28 02:41:10 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	477901	20.000	ng/u1	0.00
20) Naphthalene-d8	10.657	136	1959567	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.492	164	1176082	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.245	188	2159789	20.000	ng/u1	0.00
79) Chrysene-d12	21.392	240	1573901	20.000	ng/u1	0.00
88) Perylene-d12	23.939	264	1835334	20.000	ng/u1	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	47108	3.692	ng/uL	0.00
4) Pyridine-d5	3.687	84	260050	7.416	ng/u1	0.00
7) Phenol-d5	7.022	99	232522	5.324	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	693895	26.569	ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	752703	20.526	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	439323	12.464	ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	490563	30.821	ng/u1	0.00
24) 2-Nitrophenol-d4	9.739	143	471861	32.101	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	893044	25.182	ng/u1	0.00
31) 4-Chloroaniline-d4	10.792	131	892712	18.664	ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	3196458	30.545	ng/u1	0.00
49) Acenaphthylene-d8	14.186	160	3441179	28.859	ng/u1	0.00
54) 4-Nitrophenol-d4	14.686	143	48471	3.355	ng/u1	0.00
60) Fluorene-d10	15.486	176	2626335	30.005	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	270240	29.393	ng/u1	0.00
73) Anthracene-d10	17.345	188	3761511	32.584	ng/u1	0.00
81) Pyrene-d10	19.592	212	3903678	31.164	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.774	264	3824370	35.942	ng/u1	0.03

Target Compounds Qvalue

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

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