

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112223\
 Data File : BP018312.D
 Acq On : 23 Nov 2023 12:44
 Operator : MA/JU
 Sample : 05411-02
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0QF4

Quant Time: Nov 24 00:05:19 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 22 14:09:55 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	353381	20.000	ng/u1	0.00
20) Naphthalene-d8	10.675	136	1478351	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.504	164	938956	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	1836205	20.000	ng/u1	0.00
79) Chrysene-d12	21.357	240	1082548	20.000	ng/u1	0.00
88) Perylene-d12	23.780	264	1194723	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	37138	3.936	ng/uL	0.00
4) Pyridine-d5	3.693	84	140115	5.404	ng/u1	0.00
7) Phenol-d5	7.028	99	177836	5.507	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	570198	29.525	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.399	132	548598	20.232	ng/u1	0.00
15) 4-Methylphenol-d8	8.575	113	336727	12.920	ng/u1	-0.01
21) Nitrobenzene-d5	9.034	128	368961	30.727	ng/u1	0.00
24) 2-Nitrophenol-d4	9.752	143	297894	26.863	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.287	165	627458	23.452	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.804	131	692861	19.201	ng/u1	-0.01
46) Dimethylphthalate-d6	13.922	166	2577475	30.850	ng/u1	-0.01
49) Acenaphthylene-d8	14.198	160	2817886	29.600	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	39523	3.427	ng/u1	-0.01
60) Fluorene-d10	15.498	176	2148590	30.746	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	169556	21.692	ng/u1	-0.01
73) Anthracene-d10	17.351	188	3160979	32.208	ng/u1	-0.01
81) Pyrene-d10	19.592	212	3021282	35.067	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.621	264	2275690	32.855	ng/u1	0.00

Target Compounds Qvalue

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

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