

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112223\
 Data File : BP018311.D
 Acq On : 23 Nov 2023 12:10
 Operator : MA/JU
 Sample : 05411-01
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0QF0

Quant Time: Nov 23 23:51:12 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	351850	20.000	ng/u1	0.00
20) Naphthalene-d8	10.675	136	1502379	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.504	164	998336	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.251	188	2098874	20.000	ng/u1	-0.01
79) Chrysene-d12	21.357	240	1348631	20.000	ng/u1	0.00
88) Perylene-d12	23.780	264	1257783	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	42829	4.559	ng/uL	0.00
4) Pyridine-d5	3.687	84	223503	8.657	ng/u1	0.00
7) Phenol-d5	7.028	99	184538	5.739	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	626018	32.557	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.399	132	589860	21.848	ng/u1	0.00
15) 4-Methylphenol-d8	8.575	113	367400	14.158	ng/u1	-0.01
21) Nitrobenzene-d5	9.034	128	406053	33.275	ng/u1	0.00
24) 2-Nitrophenol-d4	9.752	143	337541	29.951	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.287	165	698179	25.678	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.810	131	973374	26.544	ng/u1	0.00
46) Dimethylphthalate-d6	13.922	166	3089039	34.774	ng/u1	-0.01
49) Acenaphthylene-d8	14.198	160	3243685	32.046	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	45495	3.710	ng/u1	-0.01
60) Fluorene-d10	15.498	176	2676594	36.024	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.610	200	233856	26.174	ng/u1	0.00
73) Anthracene-d10	17.351	188	4467295	39.821	ng/u1	-0.01
81) Pyrene-d10	19.592	212	4598671	42.844	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.621	264	3032130	41.581	ng/u1	0.00

Target Compounds Qvalue

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

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