

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
 Data File : BP018151.D
 Acq On : 14 Nov 2023 12:13
 Operator : MA/JU
 Sample : PB156906TB 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLEB906

Quant Time: Nov 15 00:37:30 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 15:50:18 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	71240	20.000	ng/u1	0.00
20) Naphthalene-d8	10.604	136	297597	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.469	164	154568	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.251	188	403701	20.000	ng/u1	0.00
79) Chrysene-d12	21.380	240	357122	20.000	ng/u1	-0.01
88) Perylene-d12	23.845	264	400602	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.158	96	14309	7.085	ng/uL	0.00
4) Pyridine-d5	3.593	84	126056	23.940	ng/u1	-0.01
7) Phenol-d5	6.952	99	210671	30.631	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	144790	35.823	ng/u1	0.00
11) 2-Chlorophenol-d4	7.310	132	169775	31.172	ng/u1	0.00
15) 4-Methylphenol-d8	8.510	113	168887	29.967	ng/u1	0.00
21) Nitrobenzene-d5	8.999	128	84647	31.571	ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	81870	26.953	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	145321	26.843	ng/u1	0.00
31) 4-Chloroaniline-d4	10.787	131	193647	26.083	ng/u1	0.00
46) Dimethylphthalate-d6	13.892	166	477770	33.500	ng/u1	0.00
49) Acenaphthylene-d8	14.163	160	489637	31.521	ng/u1	0.00
54) 4-Nitrophenol-d4	14.739	143	36259	17.014	ng/u1	0.00
60) Fluorene-d10	15.469	176	388792	33.020	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	39697	13.602	ng/u1	-0.03
73) Anthracene-d10	17.351	188	733757	33.607	ng/u1	0.00
81) Pyrene-d10	19.604	212	838528	35.152	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.686	264	789595	33.897	ng/u1	-0.02

Target Compounds Qvalue

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

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