

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012382.D
 Acq On : 31 Oct 2022 12:04
 Operator : CG/JU
 Sample : PB148438BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK438

Quant Time: Oct 31 23:49:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 28 06:05:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	297256	20.000	ng/ul	0.00
20) Naphthalene-d8	10.599	136	1264670	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.451	164	814939	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.210	188	1815127	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240	1981346	20.000	ng/ul	0.00
88) Perylene-d12	23.698	264	1965941	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	41895	5.675	ng/uL	0.00
4) Pyridine-d5	3.658	84	540884	25.505	ng/ul	-0.01
7) Phenol-d5	6.969	99	724696	27.181	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	436821	27.445	ng/ul	0.00
11) 2-Chlorophenol-d4	7.328	132	572281	29.393	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113	568801	26.974	ng/ul	-0.01
21) Nitrobenzene-d5	8.963	128	278867	34.004	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.687	143	309224	36.327	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	524954	28.457	ng/ul	0.00
31) 4-Chloroaniline-d4	10.746	131	814487	29.653	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	1782677	31.955	ng/ul	0.00
49) Acenaphthylene-d8	14.140	160	1957180	30.700	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	308769	29.244	ng/ul	0.00
60) Fluorene-d10	15.445	176	1542402	32.292	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.581	200	262039	27.728	ng/ul	0.00
73) Anthracene-d10	17.310	188	2566735	32.819	ng/ul	0.00
81) Pyrene-d10	19.551	212	3181898	31.986	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.545	264	3173355	33.036	ng/ul	0.00

Target Compounds Qvalue

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

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