

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
 Data File : BP012349.D
 Acq On : 27 Oct 2022 18:57
 Operator : CG/JU
 Sample : PB148403BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 27 23:23:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 26 23:40:20 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	330651	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.610	136	1506979	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	1041351	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.222	188	2277909	20.000	ng/ul	0.00
79) Chrysene-d12	21.316	240	2391246	20.000	ng/ul	0.00
88) Perylene-d12	23.710	264	2179619	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	41128	5.008	ng/uL	0.00
4) Pyridine-d5	3.670	84	381100	16.155	ng/ul	0.00
7) Phenol-d5	6.981	99	773822	26.093	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.146	67	474591	26.807	ng/ul	0.00
11) 2-Chlorophenol-d4	7.340	132	612449	28.279	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	630834	26.895	ng/ul	-0.01
21) Nitrobenzene-d5	8.975	128	319402	32.684	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.699	143	349959	34.502	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	621560	28.276	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	588987	17.995	ng/ul	0.00
46) Dimethylphthalate-d6	13.875	166	2128047	29.852	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	2393524	29.381	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143	377994	28.017	ng/ul	0.00
60) Fluorene-d10	15.457	176	1861236	30.495	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	305534	25.762	ng/ul	0.00
73) Anthracene-d10	17.322	188	3056841	31.145	ng/ul	0.00
81) Pyrene-d10	19.563	212	3596455	29.956	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.557	264	3271223	30.716	ng/ul	0.00

Target Compounds Qvalue

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

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