

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102422\
 Data File : BP012277.D
 Acq On : 24 Oct 2022 10:30
 Operator : CG/JU
 Sample : PB148367BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK367

Quant Time: Oct 25 05:33:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 22 01:24:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	340201	20.000	ng/u1	0.00
20) Naphthalene-d8	10.622	136	1369436	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.469	164	747873	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1315882	20.000	ng/u1	0.00
79) Chrysene-d12	21.321	240	1275132	20.000	ng/u1	0.00
88) Perylene-d12	23.721	264	1365808	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.257	96	46812	5.569	ng/uL	0.00
4) Pyridine-d5	3.681	84	673617	27.809	ng/u1	0.00
7) Phenol-d5	6.993	99	838886	28.563	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.157	67	492118	27.629	ng/u1	0.00
11) 2-Chlorophenol-d4	7.351	132	657013	29.389	ng/u1	0.00
15) 4-Methylphenol-d8	8.534	113	636220	27.872	ng/u1	0.00
21) Nitrobenzene-d5	8.987	128	319559	31.525	ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	331354	30.298	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	571575	27.862	ng/u1	0.00
31) 4-Chloroaniline-d4	10.769	131	827987	28.408	ng/u1	0.00
46) Dimethylphthalate-d6	13.886	166	1523024	29.276	ng/u1	0.00
49) Acenaphthylene-d8	14.163	160	1790639	30.155	ng/u1	0.00
54) 4-Nitrophenol-d4	14.692	143	173439	17.795	ng/u1	0.00
60) Fluorene-d10	15.469	176	1257416	28.095	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	86305	11.529	ng/u1	0.00
73) Anthracene-d10	17.327	188	1781145	30.880	ng/u1	0.00
81) Pyrene-d10	19.568	212	1970107	27.136	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.568	264	2144990	31.317	ng/u1	0.00

Target Compounds Qvalue

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

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