

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
 Data File : BP012761.D
 Acq On : 26 Nov 2022 11:20
 Operator : CG/JU
 Sample : PB149108BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK108

Quant Time: Nov 27 21:50:20 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 25 15:59:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.022	152	621076	20.000	ng/ul	0.00
20) Naphthalene-d8	10.840	136	2745866	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.657	164	1872184	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.410	188	3917583	20.000	ng/ul	0.00
79) Chrysene-d12	21.492	240	2916517	20.000	ng/ul	-0.01
88) Perylene-d12	24.010	264	2218533	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	83561	5.765	ng/uL	0.00
4) Pyridine-d5	3.822	84	1123238	25.897	ng/ul	0.00
7) Phenol-d5	7.163	99	1375070	26.943	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.346	67	881220	28.022	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	1114125	28.649	ng/ul	0.00
15) 4-Methylphenol-d8	8.716	113	1123084	27.313	ng/ul	-0.01
21) Nitrobenzene-d5	9.187	128	529131	31.484	ng/ul	0.00
24) 2-Nitrophenol-d4	9.910	143	591209	33.130	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.446	165	1136159	27.287	ng/ul	0.00
31) 4-Chloroaniline-d4	10.969	131	1632222	28.627	ng/ul	0.00
46) Dimethylphthalate-d6	14.063	166	3857547	29.935	ng/ul	0.00
49) Acenaphthylene-d8	14.351	160	4283641	28.489	ng/ul	0.00
54) 4-Nitrophenol-d4	14.834	143	528169	22.989	ng/ul	0.00
60) Fluorene-d10	15.651	176	3319158	29.410	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	424988	22.697	ng/ul	0.00
73) Anthracene-d10	17.510	188	5181885	30.049	ng/ul	0.00
81) Pyrene-d10	19.733	212	5582570	32.942	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.845	264	3337379	30.626	ng/ul	-0.01

Target Compounds Qvalue

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

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