

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
 Data File : BP012949.D
 Acq On : 05 Dec 2022 12:56
 Operator : CG/JU
 Sample : PB149301BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK301

Quant Time: Dec 05 21:50:53 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 01 23:46:27 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.005	152	537827	20.000	ng/ul	0.00
20) Naphthalene-d8	10.816	136	2535902	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.639	164	1851633	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.398	188	4072358	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	3246830	20.000	ng/ul	0.00
88) Perylene-d12	23.992	264	2924398	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	62726	5.190	ng/uL	0.00
4) Pyridine-d5	3.817	84	875963	22.458	ng/ul	0.00
7) Phenol-d5	7.158	99	1057512	21.292	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	675134	24.009	ng/ul	0.00
11) 2-Chlorophenol-d4	7.528	132	855542	23.502	ng/ul	0.00
15) 4-Methylphenol-d8	8.704	113	882270	21.634	ng/ul	0.00
21) Nitrobenzene-d5	9.169	128	439235	24.356	ng/ul	0.00
24) 2-Nitrophenol-d4	9.893	143	493470	23.911	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.434	165	897903	21.733	ng/ul	0.00
31) 4-Chloroaniline-d4	10.951	131	1311541	22.468	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	3361255	25.169	ng/ul	0.00
49) Acenaphthylene-d8	14.334	160	3599283	24.410	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	465071	17.025	ng/ul	0.00
60) Fluorene-d10	15.634	176	2864450	24.403	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	420386	17.595	ng/ul	0.00
73) Anthracene-d10	17.492	188	4696653	26.092	ng/ul	0.00
81) Pyrene-d10	19.722	212	5079240	31.826	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.827	264	3724496	26.066	ng/ul	0.00

Target Compounds Qvalue

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

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