

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120222\
 Data File : BP012930.D
 Acq On : 02 Dec 2022 09:12
 Operator : CG/JU
 Sample : PB149301BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK301

Quant Time: Dec 03 01:28:19 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	328937	20.000	ng/ul	0.00
20) Naphthalene-d8	10.822	136	1661230	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.640	164	1298610	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	3196861	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	3602602	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	3859808	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	38383	5.192	ng/uL	0.00
4) Pyridine-d5	3.817	84	524504	21.987	ng/ul	0.00
7) Phenol-d5	7.158	99	692388	22.793	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	441399	25.666	ng/ul	0.00
11) 2-Chlorophenol-d4	7.534	132	536731	24.107	ng/ul	0.00
15) 4-Methylphenol-d8	8.705	113	583401	23.390	ng/ul	0.00
21) Nitrobenzene-d5	9.169	128	280305	23.727	ng/ul	0.00
24) 2-Nitrophenol-d4	9.893	143	313224	23.168	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.434	165	606725	22.417	ng/ul	0.00
31) 4-Chloroaniline-d4	10.951	131	910584	23.812	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	2439847	26.049	ng/ul	0.00
49) Acenaphthylene-d8	14.334	160	2505050	24.224	ng/ul	0.00
54) 4-Nitrophenol-d4	14.822	143	417816	21.809	ng/ul	0.00
60) Fluorene-d10	15.634	176	2134674	25.930	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	360746	19.233	ng/ul	0.00
73) Anthracene-d10	17.492	188	3740310	26.470	ng/ul	0.00
81) Pyrene-d10	19.722	212	4628882	26.140	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	4977433	26.393	ng/ul	0.00

Target Compounds Qvalue

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

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