

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120222\
 Data File : BP012931.D
 Acq On : 02 Dec 2022 09:46
 Operator : CG/JU
 Sample : PB149337BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK337

Quant Time: Dec 03 01:28:29 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	298130	20.000	ng/u1	0.00
20) Naphthalene-d8	10.822	136	1476795	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.640	164	1134282	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.392	188	2743634	20.000	ng/u1	0.00
79) Chrysene-d12	21.475	240	3093709	20.000	ng/u1	0.00
88) Perylene-d12	23.980	264	3333776	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	53209	7.942	ng/uL	0.00
4) Pyridine-d5	3.817	84	772390	35.724	ng/u1	0.00
7) Phenol-d5	7.158	99	1014359	36.843	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	645063	41.384	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	785268	38.915	ng/u1	0.00
15) 4-Methylphenol-d8	8.711	113	861100	38.091	ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	415126	39.528	ng/u1	0.00
24) 2-Nitrophenol-d4	9.899	143	467847	38.927	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.434	165	887695	36.895	ng/u1	0.00
31) 4-Chloroaniline-d4	10.952	131	1299568	38.229	ng/u1	0.00
46) Dimethylphthalate-d6	14.046	166	3483798	42.584	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	3620410	40.081	ng/u1	0.00
54) 4-Nitrophenol-d4	14.822	143	642881	38.419	ng/u1	0.00
60) Fluorene-d10	15.634	176	3035953	42.221	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.740	200	572235	35.549	ng/u1	0.00
73) Anthracene-d10	17.492	188	5396620	44.500	ng/u1	0.00
81) Pyrene-d10	19.722	212	6730493	44.259	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.822	264	7404990	45.460	ng/u1	0.00

Target Compounds Qvalue

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

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