

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012062.D
 Acq On : 12 Oct 2022 04:17
 Operator : CG/JU
 Sample : PB148150BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK150

Quant Time: Oct 12 04:59:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 11 23:54:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	239388	20.000	ng/u1	0.00
20) Naphthalene-d8	10.663	136	962195	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.504	164	529462	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	1058707	20.000	ng/u1	-0.01
79) Chrysene-d12	21.345	240	1168505	20.000	ng/u1	0.00
88) Perylene-d12	23.768	264	1276674	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	34506	5.683	ng/uL	0.00
4) Pyridine-d5	3.699	84	425346	26.342	ng/u1	0.00
7) Phenol-d5	7.022	99	573510	29.788	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	326077	29.838	ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	477740	31.261	ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	436517	29.518	ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	215166	31.663	ng/u1	0.00
24) 2-Nitrophenol-d4	9.745	143	225625	32.744	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	395154	29.590	ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	580470	29.497	ng/u1	0.00
46) Dimethylphthalate-d6	13.916	166	1050772	31.337	ng/u1	0.00
49) Acenaphthylene-d8	14.198	160	1277860	30.818	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	184126	27.970	ng/u1	0.00
60) Fluorene-d10	15.498	176	961920	31.707	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	135354	24.212	ng/u1	0.00
73) Anthracene-d10	17.357	188	1535293	33.395	ng/u1	0.00
81) Pyrene-d10	19.598	212	1868278	31.209	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.609	264	2120969	34.175	ng/u1	0.00

Target Compounds Qvalue

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

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