

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012091.D
 Acq On : 12 Oct 2022 21:32
 Operator : CG/JU
 Sample : PB148213BL
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK213

Quant Time: Oct 12 23:10:43 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 23:03:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	246751	20.000	ng/u1	0.00
20) Naphthalene-d8	10.657	136	982822	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.498	164	554902	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	1148620	20.000	ng/u1	0.00
79) Chrysene-d12	21.345	240	1193867	20.000	ng/u1	0.00
88) Perylene-d12	23.763	264	1252979	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	39730	6.348	ng/uL	0.00
4) Pyridine-d5	3.699	84	501168	30.111	ng/u1	0.00
7) Phenol-d5	7.022	99	664365	33.477	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	374387	33.236	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	556801	35.347	ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	512627	33.630	ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	254551	36.673	ng/u1	0.00
24) 2-Nitrophenol-d4	9.746	143	276843	39.334	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	468306	34.332	ng/u1	0.00
31) 4-Chloroaniline-d4	10.805	131	713335	35.488	ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	1289670	36.698	ng/u1	0.00
49) Acenaphthylene-d8	14.193	160	1578941	36.334	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	229425	33.253	ng/u1	0.00
60) Fluorene-d10	15.498	176	1189723	37.418	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	166507	27.453	ng/u1	0.00
73) Anthracene-d10	17.357	188	1895704	38.007	ng/u1	0.00
81) Pyrene-d10	19.592	212	2262998	37.000	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.610	264	2345269	38.504	ng/u1	0.00

Target Compounds Qvalue

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

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