

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112922\
 Data File : BP012858.D
 Acq On : 29 Nov 2022 18:37
 Operator : CG/JU
 Sample : PB149234BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK234

Quant Time: Nov 29 22:58:14 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 29 22:53:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	571258	20.000	ng/u1	0.00
20) Naphthalene-d8	10.828	136	2609179	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.645	164	1786135	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	3809310	20.000	ng/u1	0.00
79) Chrysene-d12	21.486	240	2923991	20.000	ng/u1	0.00
88) Perylene-d12	23.998	264	2358582	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	80511	6.039	ng/uL	0.00
4) Pyridine-d5	3.817	84	762656	19.117	ng/u1	0.00
7) Phenol-d5	7.163	99	1401533	29.857	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	906969	31.355	ng/u1	0.00
11) 2-Chlorophenol-d4	7.540	132	1127666	31.526	ng/u1	0.00
15) 4-Methylphenol-d8	8.710	113	1150238	30.412	ng/u1	0.00
21) Nitrobenzene-d5	9.181	128	551617	34.542	ng/u1	0.00
24) 2-Nitrophenol-d4	9.905	143	601247	35.458	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	1166022	29.471	ng/u1	0.00
31) 4-Chloroaniline-d4	10.957	131	1381089	25.491	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	4021875	32.713	ng/u1	0.00
49) Acenaphthylene-d8	14.340	160	4443745	30.978	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	572146	26.103	ng/u1	0.00
60) Fluorene-d10	15.640	176	3502640	32.531	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	483605	26.562	ng/u1	0.00
73) Anthracene-d10	17.504	188	5497126	32.783	ng/u1	0.00
81) Pyrene-d10	19.728	212	5858535	34.482	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.833	264	3776545	32.598	ng/u1	0.00

Target Compounds Qvalue

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

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