

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012109.D
 Acq On : 13 Oct 2022 08:24
 Operator : CG/JU
 Sample : PB148251BL
 Misc :
 ALS Vial : 80 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK251

Quant Time: Oct 13 08:51:00 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	7.846	152	223547	20.000	ng/u1	0.00
20) Naphthalene-d8	10.651	136	902993	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.498	164	519458	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	1109745	20.000	ng/u1	0.00
79) Chrysene-d12	21.345	240	1159881	20.000	ng/u1	0.00
88) Perylene-d12	23.763	264	1215101	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	37551	6.623	ng/uL	0.00
4) Pyridine-d5	3.699	84	464710	30.819	ng/u1	0.00
7) Phenol-d5	7.016	99	618815	34.419	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	347584	34.060	ng/u1	0.00
11) 2-Chlorophenol-d4	7.375	132	510473	35.769	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	468708	33.941	ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	236800	37.131	ng/u1	0.00
24) 2-Nitrophenol-d4	9.740	143	251180	38.843	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	439121	35.039	ng/u1	0.00
31) 4-Chloroaniline-d4	10.799	131	654277	35.427	ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	1196911	36.382	ng/u1	0.00
49) Acenaphthylene-d8	14.192	160	1449263	35.625	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	216999	33.598	ng/u1	0.00
60) Fluorene-d10	15.492	176	1122702	37.719	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.616	200	154442	26.356	ng/u1	0.00
73) Anthracene-d10	17.357	188	1821935	37.808	ng/u1	0.00
81) Pyrene-d10	19.592	212	2178621	36.664	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.604	264	2257865	38.224	ng/u1	0.00

Target Compounds Qvalue

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

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