

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082222\
 Data File : BP011531.D
 Acq On : 22 Aug 2022 14:14
 Operator : CG/JU
 Sample : PB146658BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK658

Quant Time: Aug 22 15:14:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 17 01:58:37 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.552	152	54707	20.000 ng/u1	0.00
20) Naphthalene-d8	10.334	136	232405	20.000 ng/u1	-0.02
38) Acenaphthene-d10	14.222	164	134923	20.000 ng/u1	-0.01
64) Phenanthrene-d10	16.992	188	268053	20.000 ng/u1	-0.01
79) Chrysene-d12	21.121	240	242037	20.000 ng/u1	-0.01
88) Perylene-d12	23.410	264	249643	20.000 ng/u1	-0.02
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.064	96	10635	6.033 ng/uL	0.00
4) Pyridine-d5	3.475	84	111818	24.774 ng/u1	0.00
7) Phenol-d5	6.734	99	150731	29.982 ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	6.899	67	101416	31.050 ng/u1	-0.01
11) 2-Chlorophenol-d4	7.081	132	125397	33.913 ng/u1	-0.01
15) 4-Methylphenol-d8	8.269	113	123545	30.703 ng/u1	-0.01
21) Nitrobenzene-d5	8.716	128	59394	35.498 ng/u1	-0.01
24) 2-Nitrophenol-d4	9.428	143	61863	36.330 ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	9.963	165	104236	33.162 ng/u1	-0.01
31) 4-Chloroaniline-d4	10.493	131	182598	34.898 ng/u1	0.00
46) Dimethylphthalate-d6	13.645	166	355559	37.259 ng/u1	-0.01
49) Acenaphthylene-d8	13.910	160	401224	37.829 ng/u1	-0.01
54) 4-Nitrophenol-d4	14.451	143	51317	27.911 ng/u1	-0.01
60) Fluorene-d10	15.228	176	284899	35.487 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.363	200	44928	31.018 ng/u1	0.00
73) Anthracene-d10	17.092	188	447948	37.732 ng/u1	-0.01
81) Pyrene-d10	19.357	212	501352	40.941 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.262	264	474238	39.053 ng/u1	-0.02

Target Compounds Qvalue

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

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