

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082322\
 Data File : BP011541.D
 Acq On : 23 Aug 2022 14:37
 Operator : CG/JU
 Sample : PB146669BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK669

Quant Time: Aug 23 16:30:52 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	55836	20.000 ng/u1	0.00
20) Naphthalene-d8	10.340	136	236618	20.000 ng/u1	-0.01
38) Acenaphthene-d10	14.222	164	135699	20.000 ng/u1	-0.01
64) Phenanthrene-d10	16.998	188	275077	20.000 ng/u1	0.00
79) Chrysene-d12	21.128	240	256455	20.000 ng/u1	0.00
88) Perylene-d12	23.416	264	273338	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.064	96	9513	5.287 ng/uL	0.00
4) Pyridine-d5	3.476	84	114465	24.847 ng/u1	0.00
7) Phenol-d5	6.734	99	129323	25.204 ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	6.899	67	86794	26.036 ng/u1	-0.01
11) 2-Chlorophenol-d4	7.087	132	106025	28.094 ng/u1	0.00
15) 4-Methylphenol-d8	8.269	113	104342	25.406 ng/u1	-0.01
21) Nitrobenzene-d5	8.716	128	49799	29.234 ng/u1	-0.01
24) 2-Nitrophenol-d4	9.434	143	52329	30.184 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.969	165	87541	27.355 ng/u1	0.00
31) 4-Chloroaniline-d4	10.493	131	139154	26.122 ng/u1	0.00
46) Dimethylphthalate-d6	13.651	166	315365	32.858 ng/u1	0.00
49) Acenaphthylene-d8	13.916	160	325096	30.476 ng/u1	0.00
54) 4-Nitrophenol-d4	14.457	143	43832	23.703 ng/u1	0.00
60) Fluorene-d10	15.228	176	250233	30.990 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	40743	27.411 ng/u1	0.00
73) Anthracene-d10	17.098	188	400861	32.904 ng/u1	0.00
81) Pyrene-d10	19.363	212	453488	34.950 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.269	264	457834	34.433 ng/u1	-0.01

Target Compounds Qvalue

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

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