

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082222\
 Data File : BP011532.D
 Acq On : 22 Aug 2022 14:49
 Operator : CG/JU
 Sample : PB146659BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK659

Quant Time: Aug 22 15:17:56 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	52117	20.000 ng/u1	0.00
20) Naphthalene-d8	10.340	136	230932	20.000 ng/u1	-0.01
38) Acenaphthene-d10	14.222	164	132187	20.000 ng/u1	-0.01
64) Phenanthrene-d10	16.992	188	260552	20.000 ng/u1	-0.01
79) Chrysene-d12	21.127	240	235482	20.000 ng/u1	0.00
88) Perylene-d12	23.415	264	246729	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.069	96	10606	6.315 ng/uL	0.00
4) Pyridine-d5	3.475	84	108719	25.284 ng/u1	0.00
7) Phenol-d5	6.734	99	147012	30.696 ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	6.899	67	98731	31.731 ng/u1	-0.01
11) 2-Chlorophenol-d4	7.081	132	121999	34.634 ng/u1	-0.01
15) 4-Methylphenol-d8	8.269	113	119766	31.243 ng/u1	-0.01
21) Nitrobenzene-d5	8.716	128	57809	34.771 ng/u1	-0.01
24) 2-Nitrophenol-d4	9.434	143	61475	36.332 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.963	165	103418	33.112 ng/u1	-0.01
31) 4-Chloroaniline-d4	10.493	131	176335	33.916 ng/u1	0.00
46) Dimethylphthalate-d6	13.651	166	349280	37.359 ng/u1	0.00
49) Acenaphthylene-d8	13.910	160	396176	38.126 ng/u1	-0.01
54) 4-Nitrophenol-d4	14.457	143	46638	25.891 ng/u1	0.00
60) Fluorene-d10	15.228	176	282676	35.939 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.363	200	43628	30.988 ng/u1	0.00
73) Anthracene-d10	17.092	188	441513	38.261 ng/u1	-0.01
81) Pyrene-d10	19.357	212	484406	40.658 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.268	264	460925	38.405 ng/u1	-0.01

Target Compounds Qvalue

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

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