

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
 Data File : BP011515.D
 Acq On : 19 Aug 2022 20:06
 Operator : CG/JU
 Sample : PB147110BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK110

Quant Time: Aug 20 04:40:13 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	105182	20.000 ng/u1	0.00
20) Naphthalene-d8	10.340	136	437306	20.000 ng/u1	-0.01
38) Acenaphthene-d10	14.228	164	248161	20.000 ng/u1	0.00
64) Phenanthrene-d10	16.998	188	493853	20.000 ng/u1	0.00
79) Chrysene-d12	21.127	240	467054	20.000 ng/u1	0.00
88) Perylene-d12	23.415	264	484689	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.069	96	15514	4.577 ng/uL	0.00
4) Pyridine-d5	3.481	84	190605	21.964 ng/u1	0.00
7) Phenol-d5	6.740	99	223898	23.164 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.905	67	148914	23.714 ng/u1	0.00
11) 2-Chlorophenol-d4	7.087	132	180406	25.377 ng/u1	0.00
15) 4-Methylphenol-d8	8.275	113	179284	23.174 ng/u1	0.00
21) Nitrobenzene-d5	8.716	128	85528	27.166 ng/u1	-0.01
24) 2-Nitrophenol-d4	9.434	143	88399	27.589 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.969	165	148852	25.168 ng/u1	0.00
31) 4-Chloroaniline-d4	10.493	131	232999	23.666 ng/u1	0.00
46) Dimethylphthalate-d6	13.651	166	499128	28.437 ng/u1	0.00
49) Acenaphthylene-d8	13.916	160	525914	26.959 ng/u1	0.00
54) 4-Nitrophenol-d4	14.457	143	77312	22.862 ng/u1	0.00
60) Fluorene-d10	15.228	176	387756	26.259 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	65633	24.595 ng/u1	0.00
73) Anthracene-d10	17.098	188	633430	28.961 ng/u1	0.00
81) Pyrene-d10	19.363	212	713247	30.183 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.268	264	713868	30.278 ng/u1	-0.01

Target Compounds Qvalue

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

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