

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081722\
 Data File : BP011454.D
 Acq On : 17 Aug 2022 10:13
 Operator : CG/JU
 Sample : PB147024BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Aug 17 23:24:32 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	7.558	152	77269	20.000	ng/u1	0.00
20) Naphthalene-d8	10.346	136	338229	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.228	164	203237	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.998	188	409023	20.000	ng/u1	0.00
79) Chrysene-d12	21.133	240	377919	20.000	ng/u1	0.00
88) Perylene-d12	23.421	264	385846	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.070	96	14211	5.707	ng/uL	0.00
4) Pyridine-d5	3.475	84	175029	27.455	ng/u1	0.00
7) Phenol-d5	6.740	99	197423	27.803	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.905	67	133236	28.881	ng/u1	0.00
11) 2-Chlorophenol-d4	7.087	132	158814	30.409	ng/u1	0.00
15) 4-Methylphenol-d8	8.275	113	162802	28.645	ng/u1	0.00
21) Nitrobenzene-d5	8.722	128	76526	31.427	ng/u1	0.00
24) 2-Nitrophenol-d4	9.440	143	79186	31.953	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	133928	29.278	ng/u1	0.00
31) 4-Chloroaniline-d4	10.499	131	226782	29.782	ng/u1	0.00
46) Dimethylphthalate-d6	13.657	166	528433	36.761	ng/u1	0.00
49) Acenaphthylene-d8	13.922	160	555252	34.755	ng/u1	0.00
54) 4-Nitrophenol-d4	14.457	143	73678	26.603	ng/u1	0.00
60) Fluorene-d10	15.234	176	411596	34.035	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	66319	30.006	ng/u1	0.00
73) Anthracene-d10	17.104	188	663288	36.615	ng/u1	0.00
81) Pyrene-d10	19.363	212	745567	38.993	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.274	264	704910	37.557	ng/u1	0.00

Target Compounds	Qvalue

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

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