

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080522\
 Data File : BP011274.D
 Acq On : 05 Aug 2022 11:58
 Operator : CG/JU
 Sample : PB146743BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Aug 05 23:01:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080222.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 02 23:52:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.598	152	108610	20.000	ng/u1	0.00
20) Naphthalene-d8	10.392	136	440501	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.275	164	240462	20.000	ng/u1	# 0.00
64) Phenanthrene-d10	17.045	188	485049	20.000	ng/u1	0.00
79) Chrysene-d12	21.168	240	644210	20.000	ng/u1	0.00
88) Perylene-d12	23.486	264	695265	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	5673	4.223	ng/uL	0.00
4) Pyridine-d5	3.505	84	66182	18.266	ng/u1	0.00
7) Phenol-d5	6.781	99	237764	26.269	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67	233107	35.691	ng/u1	0.00
11) 2-Chlorophenol-d4	7.134	132	199894	27.095	ng/u1	0.00
15) 4-Methylphenol-d8	8.322	113	189827	25.489	ng/u1	0.00
21) Nitrobenzene-d5	8.763	128	143374	45.251	ng/u1	0.00
24) 2-Nitrophenol-d4	9.481	143	95985	32.319	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	160284	28.301	ng/u1	0.00
31) 4-Chloroaniline-d4	10.545	131	235713	26.903	ng/u1	0.00
46) Dimethylphthalate-d6	13.698	166	531727	32.286	ng/u1	0.00
49) Acenaphthylene-d8	13.963	160	827913	39.955	ng/u1	0.00
54) 4-Nitrophenol-d4	14.504	143	76346	27.261	ng/u1	0.00
60) Fluorene-d10	15.275	176	678069	48.387	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	71017	31.531	ng/u1	0.00
73) Anthracene-d10	17.139	188	1262195	59.209	ng/u1	0.00
81) Pyrene-d10	19.404	212	1412722	43.373	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.339	264	1505274	43.447	ng/u1	0.00

Target Compounds Qvalue

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

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