

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
 Data File : BP010474.D
 Acq On : 23 May 2022 09:59
 Operator : CG/JU
 Sample : PB144961BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: May 23 11:07:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.892	152	135959	20.000	ng/u1	0.00
20) Naphthalene-d8	10.692	136	613145	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.521	164	412207	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.274	188	913390	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.356	240	865871	20.000	ng/u1	# 0.00
88) Perylene-d12	23.786	264	753618	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	18391	5.766	ng/uL	0.00
4) Pyridine-d5	3.757	84	267669	29.555	ng/u1	0.00
7) Phenol-d5	7.057	99	363394	31.845	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	245107	31.969	ng/u1	0.00
11) 2-Chlorophenol-d4	7.422	132	292686	33.519	ng/u1	-0.01
15) 4-Methylphenol-d8	8.604	113	315968	32.581	ng/u1	0.00
21) Nitrobenzene-d5	9.051	128	157904	33.276	ng/u1	0.00
24) 2-Nitrophenol-d4	9.775	143	170843	33.972	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	295300	31.055	ng/u1	0.00
31) 4-Chloroaniline-d4	10.828	131	431091	30.490	ng/u1	-0.01
46) Dimethylphthalate-d6	13.933	166	1022997	33.684	ng/u1	0.00
49) Acenaphthylene-d8	14.216	160	1188234m	33.919	ng/u1	0.00
54) 4-Nitrophenol-d4	14.727	143	134141	24.815	ng/u1	0.00
60) Fluorene-d10	15.516	176	880988	33.344	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	130884	24.502	ng/u1	0.00
73) Anthracene-d10	17.374	188	1424775	34.448	ng/u1	0.00
81) Pyrene-d10	19.609	212	1753274	38.128	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.627	264	1372920	36.280	ng/u1	0.00

Target Compounds	Qvalue

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

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