

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052022\
 Data File : BP010458.D
 Acq On : 21 May 2022 04:45
 Operator : CG/JU
 Sample : PB144914BL
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK914

Quant Time: May 21 06:04:56 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	7.893	152	167445	20.000	ng/u1	0.00
20) Naphthalene-d8	10.693	136	756578	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.522	164	515967	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.275	188	1124739	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.357	240	1102690	20.000	ng/u1	# 0.00
88) Perylene-d12	23.780	264	982462	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	22707	5.780	ng/uL	0.00
4) Pyridine-d5	3.758	84	333281	29.880	ng/u1	0.00
7) Phenol-d5	7.052	99	416761	29.654	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.216	67	291199	30.839	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.422	132	335877	31.233	ng/u1	-0.01
15) 4-Methylphenol-d8	8.599	113	357849	29.961	ng/u1	-0.01
21) Nitrobenzene-d5	9.046	128	182977	31.249	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.775	143	198311	31.958	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	340448	29.015	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.828	131	503365	28.853	ng/u1	-0.01
46) Dimethylphthalate-d6	13.928	166	1224403	32.209	ng/u1	-0.01
49) Acenaphthylene-d8	14.216	160	1381773	31.512	ng/u1	0.00
54) 4-Nitrophenol-d4	14.716	143	172279	25.461	ng/u1	-0.01
60) Fluorene-d10	15.516	176	1041626	31.496	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	162012	24.630	ng/u1	-0.01
73) Anthracene-d10	17.375	188	1651758	32.432	ng/u1	0.00
81) Pyrene-d10	19.604	212	2034706	34.745	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.621	264	1633330	33.108	ng/u1	-0.01

Target Compounds Qvalue

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

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