

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
 Data File : BP010922.D
 Acq On : 30 Jun 2022 17:27
 Operator : CG/JU
 Sample : PB145664BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK664

Quant Time: Jul 01 02:33:47 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 27 23:12:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	386205	20.000	ng/u1	0.00
20) Naphthalene-d8	10.499	136	1546507	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.363	164	769223	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.122	188	1380479	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.239	240	1389101	20.000	ng/u1	# 0.00
88) Perylene-d12	23.598	264	1525123	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	75267	7.372	ng/uL	0.00
4) Pyridine-d5	3.605	84	971759	35.757	ng/u1	0.00
7) Phenol-d5	6.881	99	1140452	37.303	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	729803	36.501	ng/u1	0.00
11) 2-Chlorophenol-d4	7.240	132	961239	37.897	ng/u1	0.00
15) 4-Methylphenol-d8	8.416	113	885381	36.635	ng/u1	-0.01
21) Nitrobenzene-d5	8.863	128	426474	43.319	ng/u1	0.00
24) 2-Nitrophenol-d4	9.587	143	409224	50.242	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.122	165	722297	35.890	ng/u1	0.00
31) 4-Chloroaniline-d4	10.646	131	1176378	35.809	ng/u1	0.00
46) Dimethylphthalate-d6	13.781	166	1929614	36.620	ng/u1	0.00
49) Acenaphthylene-d8	14.051	160	2424300	36.049	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	305421	37.110	ng/u1	0.00
60) Fluorene-d10	15.357	176	1666777	36.861	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.481	200	144847	31.670	ng/u1	0.00
73) Anthracene-d10	17.222	188	2359467	38.416	ng/u1	-0.01
81) Pyrene-d10	19.475	212	2645041	34.951	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.445	264	3024934	40.137	ng/u1	-0.01

Target Compounds	Qvalue

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

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