

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
 Data File : BP010912.D
 Acq On : 30 Jun 2022 10:22
 Operator : CG/JU
 Sample : PB145923BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK923

Quant Time: Jul 01 02:31:53 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	469789	20.000	ng/u1	0.00
20) Naphthalene-d8	10.499	136	1876763	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.363	164	949515	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.128	188	1694682	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.239	240	1693051	20.000	ng/u1	# 0.00
88) Perylene-d12	23.598	264	1855227	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	70404	5.669	ng/uL	0.00
4) Pyridine-d5	3.611	84	900948	27.253	ng/u1	0.00
7) Phenol-d5	6.881	99	1037066	27.886	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	666210	27.392	ng/u1	0.00
11) 2-Chlorophenol-d4	7.240	132	869748	28.189	ng/u1	0.00
15) 4-Methylphenol-d8	8.422	113	805224	27.390	ng/u1	0.00
21) Nitrobenzene-d5	8.869	128	383506	32.100	ng/u1	0.00
24) 2-Nitrophenol-d4	9.587	143	364068	36.833	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.122	165	652852	26.731	ng/u1	0.00
31) 4-Chloroaniline-d4	10.646	131	1050850	26.359	ng/u1	0.00
46) Dimethylphthalate-d6	13.781	166	1767377	27.173	ng/u1	0.00
49) Acenaphthylene-d8	14.051	160	2141593	25.799	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	273399	26.911	ng/u1	0.00
60) Fluorene-d10	15.357	176	1472719	26.385	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.481	200	124528	22.179	ng/u1	0.00
73) Anthracene-d10	17.222	188	2211563	29.332	ng/u1	-0.01
81) Pyrene-d10	19.475	212	2507272	27.183	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.445	264	2812531	30.679	ng/u1	-0.01

Target Compounds Qvalue

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

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