

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010329.D
 Acq On : 12 May 2022 18:34
 Operator : CG/JU
 Sample : PB144737BL
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK737

Quant Time: May 13 00:02:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 12 01:10:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	142181	20.000	ng/u1	# 0.00
20) Naphthalene-d8	10.705	136	633220	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.534	164	451648	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.287	188	998007	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.363	240	1025063	20.000	ng/u1	#-0.01
88) Perylene-d12	23.792	264	989021	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	25379	7.583	ng/uL	0.00
4) Pyridine-d5	3.764	84	374908	40.254	ng/u1	0.00
7) Phenol-d5	7.069	99	481708	36.478	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	339553	41.564	ng/u1	0.00
11) 2-Chlorophenol-d4	7.434	132	388116	39.907	ng/u1	0.00
15) 4-Methylphenol-d8	8.611	113	422946	38.957	ng/u1	0.00
21) Nitrobenzene-d5	9.063	128	211790	41.713	ng/u1	0.00
24) 2-Nitrophenol-d4	9.787	143	234093	40.438	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	394786	37.468	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.840	131	600866	38.628	ng/u1	0.00
46) Dimethylphthalate-d6	13.946	166	1517456	42.173	ng/u1	0.00
49) Acenaphthylene-d8	14.228	160	1656714	38.619	ng/u1	0.00
54) 4-Nitrophenol-d4	14.728	143	241808	35.036	ng/u1	-0.01
60) Fluorene-d10	15.528	176	1265621	41.670	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	239335	35.020	ng/u1	-0.01
73) Anthracene-d10	17.387	188	2069534	42.345	ng/u1	0.00
81) Pyrene-d10	19.616	212	2589210	45.709	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.639	264	2439051	46.243	ng/u1	-0.02

Target Compounds	Qvalue

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

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