

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
 Data File : BP010997.D
 Acq On : 18 Jul 2022 17:35
 Operator : CG/JU
 Sample : PB146279BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK279

Quant Time: Jul 19 00:46:14 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 14 15:34:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.681	152	311713	20.000	ng/u1	0.00
20) Naphthalene-d8	10.475	136	1351921	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.339	164	752329	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.110	188	1534837	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.227	240	1562169	20.000	ng/u1	#-0.01
88) Perylene-d12	23.580	264	1653710	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	56430	6.526	ng/uL	0.00
4) Pyridine-d5	3.587	84	652909	29.215	ng/u1	0.00
7) Phenol-d5	6.858	99	774610	30.273	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.022	67	528758	32.136	ng/u1	0.00
11) 2-Chlorophenol-d4	7.216	132	658510	31.850	ng/u1	0.00
15) 4-Methylphenol-d8	8.393	113	602176	29.397	ng/u1	-0.01
21) Nitrobenzene-d5	8.840	128	291480	29.585	ng/u1	0.00
24) 2-Nitrophenol-d4	9.563	143	271231	27.822	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.099	165	478810	26.857	ng/u1	0.00
31) 4-Chloroaniline-d4	10.622	131	654674	22.274	ng/u1	0.00
46) Dimethylphthalate-d6	13.763	166	1611280	31.336	ng/u1	0.00
49) Acenaphthylene-d8	14.034	160	1865378	30.747	ng/u1	0.00
54) 4-Nitrophenol-d4	14.557	143	252030	24.218	ng/u1	0.00
60) Fluorene-d10	15.339	176	1345462	30.784	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.463	200	170456	22.444	ng/u1	-0.01
73) Anthracene-d10	17.210	188	2125087	32.569	ng/u1	0.00
81) Pyrene-d10	19.463	212	2496570	30.429	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.427	264	2569263	32.348	ng/u1	-0.02

Target Compounds	Qvalue

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

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