

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
 Data File : BP010996.D
 Acq On : 18 Jul 2022 17:01
 Operator : CG/JU
 Sample : PB146261BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK261

Quant Time: Jul 19 00:46:06 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	331177	20.000	ng/u1	0.00
20) Naphthalene-d8	10.475	136	1372652	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.339	164	714901	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.110	188	1390686	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.227	240	1430716	20.000	ng/u1	#-0.01
88) Perylene-d12	23.580	264	1557963	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	61619	6.708	ng/uL	0.00
4) Pyridine-d5	3.587	84	688583	29.000	ng/u1	0.00
7) Phenol-d5	6.858	99	839087	30.866	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.022	67	570428	32.631	ng/u1	0.00
11) 2-Chlorophenol-d4	7.216	132	720491	32.800	ng/u1	0.00
15) 4-Methylphenol-d8	8.393	113	641965	29.497	ng/u1	-0.01
21) Nitrobenzene-d5	8.846	128	311023	31.092	ng/u1	0.00
24) 2-Nitrophenol-d4	9.563	143	291420	29.442	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.099	165	519664	28.708	ng/u1	0.00
31) 4-Chloroaniline-d4	10.622	131	705153	23.629	ng/u1	0.00
46) Dimethylphthalate-d6	13.763	166	1548408	31.690	ng/u1	0.00
49) Acenaphthylene-d8	14.034	160	1864767	32.346	ng/u1	0.00
54) 4-Nitrophenol-d4	14.557	143	228041	23.060	ng/u1	0.00
60) Fluorene-d10	15.339	176	1317829	31.730	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.463	200	154217	22.410	ng/u1	-0.01
73) Anthracene-d10	17.210	188	1961812	33.183	ng/u1	0.00
81) Pyrene-d10	19.463	212	2275361	30.281	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.427	264	2505136	33.479	ng/u1	-0.02

Target Compounds Qvalue

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

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