

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071122\
 Data File : BP010956.D
 Acq On : 11 Jul 2022 19:16
 Operator : CG/JU
 Sample : PB146106TB 10X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLEB106

Quant Time: Jul 12 01:05:57 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 06 23:13:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	473467	20.000	ng/u1	0.00
20) Naphthalene-d8	10.493	136	1889505	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.351	164	916667	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.116	188	1644910	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.227	240	1605381	20.000	ng/u1	#-0.01
88) Perylene-d12	23.586	264	1799044	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	80833	6.458	ng/uL	0.00
4) Pyridine-d5	3.605	84	863934	25.931	ng/u1	0.00
7) Phenol-d5	6.875	99	1151128	30.713	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67	751503	30.659	ng/u1	0.00
11) 2-Chlorophenol-d4	7.234	132	969695	31.184	ng/u1	0.00
15) 4-Methylphenol-d8	8.410	113	885748	29.895	ng/u1	-0.01
21) Nitrobenzene-d5	8.857	128	414380	34.450	ng/u1	0.00
24) 2-Nitrophenol-d4	9.581	143	380072	38.192	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.116	165	700793	28.500	ng/u1	0.00
31) 4-Chloroaniline-d4	10.634	131	831535	20.717	ng/u1	0.00
46) Dimethylphthalate-d6	13.775	166	1754569	27.942	ng/u1	0.00
49) Acenaphthylene-d8	14.045	160	2263558	28.245	ng/u1	0.00
54) 4-Nitrophenol-d4	14.563	143	277866	28.331	ng/u1	-0.01
60) Fluorene-d10	15.351	176	1563152	29.009	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	145830	26.759	ng/u1	0.00
73) Anthracene-d10	17.216	188	2257896	30.852	ng/u1	0.00
81) Pyrene-d10	19.469	212	2443485	27.938	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.433	264	2765055	31.103	ng/u1	-0.01

Target Compounds Qvalue

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

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