

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071322\
 Data File : BP010962.D
 Acq On : 13 Jul 2022 11:53
 Operator : CG/JU
 Sample : PB146170BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK170

Quant Time: Jul 14 02:44:39 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	495252	20.000	ng/u1	0.00
20) Naphthalene-d8	10.493	136	2194061	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.357	164	1304925	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.122	188	2657965	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.233	240	2508849	20.000	ng/u1	# 0.00
88) Perylene-d12	23.592	264	2426230	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	59781	4.566	ng/uL	0.00
4) Pyridine-d5	3.605	84	782186	22.444	ng/u1	0.00
7) Phenol-d5	6.875	99	946514	24.143	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67	625988	24.415	ng/u1	0.00
11) 2-Chlorophenol-d4	7.234	132	775970	23.857	ng/u1	0.00
15) 4-Methylphenol-d8	8.416	113	764114	24.656	ng/u1	0.00
21) Nitrobenzene-d5	8.863	128	361355	25.871	ng/u1	0.00
24) 2-Nitrophenol-d4	9.581	143	340944	29.505	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.116	165	610288	21.374	ng/u1	0.00
31) 4-Chloroaniline-d4	10.640	131	1081894	23.213	ng/u1	0.00
46) Dimethylphthalate-d6	13.775	166	2161401	24.180	ng/u1	0.00
49) Acenaphthylene-d8	14.045	160	2444583	21.428	ng/u1	0.00
54) 4-Nitrophenol-d4	14.569	143	346855	24.843	ng/u1	0.00
60) Fluorene-d10	15.357	176	1805170	23.533	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	224691	25.515	ng/u1	0.00
73) Anthracene-d10	17.222	188	2896051	24.490	ng/u1	0.00
81) Pyrene-d10	19.469	212	3384270	24.760	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.439	264	3199646	26.687	ng/u1	0.00

Target Compounds Qvalue

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

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