

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
 Data File : BP010889.D
 Acq On : 27 Jun 2022 22:11
 Operator : CG/JU
 Sample : PB145713BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK713

Quant Time: Jun 27 23:17:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 27 23:12:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	471276	20.000	ng/u1	0.00
20) Naphthalene-d8	10.504	136	1942450	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.363	164	1014871	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.127	188	1811790	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.239	240	1504762	20.000	ng/u1	# 0.00
88) Perylene-d12	23.604	264	1630553	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	74042	5.943	ng/uL	0.00
4) Pyridine-d5	3.611	84	977854	29.486	ng/u1	0.00
7) Phenol-d5	6.881	99	1109943	29.752	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.052	67	729354	29.894	ng/u1	0.00
11) 2-Chlorophenol-d4	7.240	132	929204	30.021	ng/u1	0.00
15) 4-Methylphenol-d8	8.422	113	877094	29.741	ng/u1	0.00
21) Nitrobenzene-d5	8.869	128	415603	33.610	ng/u1	0.00
24) 2-Nitrophenol-d4	9.587	143	381442	37.285	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.122	165	706328	27.943	ng/u1	0.00
31) 4-Chloroaniline-d4	10.645	131	1190415	28.850	ng/u1	0.00
46) Dimethylphthalate-d6	13.786	166	2025588	29.137	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	2440109	27.502	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	289973	26.705	ng/u1	0.00
60) Fluorene-d10	15.363	176	1725631	28.925	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.481	200	153778	25.618	ng/u1	0.00
73) Anthracene-d10	17.227	188	2464003	30.567	ng/u1	0.00
81) Pyrene-d10	19.474	212	2587490	31.562	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.451	264	2569297	31.887	ng/u1	0.00

Target Compounds	Qvalue

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

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