

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010839.D
 Acq On : 25 Jun 2022 21:49
 Operator : CG/JU
 Sample : PB145708TB 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLEB708

Quant Time: Jun 26 22:54:31 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 24 22:42:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	415505	20.000	ng/u1	0.00
20) Naphthalene-d8	10.510	136	1665209	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.375	164	876410	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.133	188	1628320	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.245	240	1613125	20.000	ng/u1	#-0.01
88) Perylene-d12	23.610	264	1734010	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	79064	7.198	ng/uL	0.00
4) Pyridine-d5	3.611	84	982438	33.601	ng/u1	0.00
7) Phenol-d5	6.899	99	1151817	35.019	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.057	67	733613	34.104	ng/u1	0.00
11) 2-Chlorophenol-d4	7.252	132	982373	35.999	ng/u1	0.00
15) 4-Methylphenol-d8	8.434	113	912157	35.081	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	441084	41.609	ng/u1	0.00
24) 2-Nitrophenol-d4	9.598	143	422136	48.133	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.134	165	759302	35.039	ng/u1	0.00
31) 4-Chloroaniline-d4	10.657	131	1229912	34.770	ng/u1	0.00
46) Dimethylphthalate-d6	13.792	166	2181255	36.333	ng/u1	0.00
49) Acenaphthylene-d8	14.063	160	2677589	34.946	ng/u1	0.00
54) 4-Nitrophenol-d4	14.586	143	343749	36.659	ng/u1	0.00
60) Fluorene-d10	15.369	176	1851855	35.945	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.492	200	162151	30.057	ng/u1	0.00
73) Anthracene-d10	17.233	188	2701696	37.292	ng/u1	-0.01
81) Pyrene-d10	19.486	212	3027226	34.446	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.457	264	3238216	37.791	ng/u1	-0.02

Target Compounds Qvalue

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

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