

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
 Data File : BP010888.D
 Acq On : 27 Jun 2022 21:37
 Operator : CG/JU
 Sample : PB145835BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK835

Quant Time: Jun 27 23:17:28 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	7.711	152	555546	20.000	ng/u1	0.00
20) Naphthalene-d8	10.505	136	2296109	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.363	164	1228303	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.128	188	2171175	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.239	240	1621237	20.000	ng/u1	# 0.00
88) Perylene-d12	23.604	264	1686774	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	90561	6.166	ng/uL	0.00
4) Pyridine-d5	3.611	84	1186106	30.341	ng/u1	0.00
7) Phenol-d5	6.881	99	1392337	31.660	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.052	67	905007	31.467	ng/u1	0.00
11) 2-Chlorophenol-d4	7.240	132	1165153	31.934	ng/u1	0.00
15) 4-Methylphenol-d8	8.422	113	1110452	31.942	ng/u1	0.00
21) Nitrobenzene-d5	8.869	128	531776	36.381	ng/u1	0.00
24) 2-Nitrophenol-d4	9.587	143	508895	42.082	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.128	165	921072	30.826	ng/u1	0.00
31) 4-Chloroaniline-d4	10.646	131	1518285	31.129	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2702805	32.123	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	3272386	30.473	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	384039	29.222	ng/u1	0.00
60) Fluorene-d10	15.363	176	2245423	31.098	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.481	200	210540	29.269	ng/u1	0.00
73) Anthracene-d10	17.228	188	3170199	32.818	ng/u1	0.00
81) Pyrene-d10	19.475	212	3138998	35.539	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.451	264	2841177	34.086	ng/u1	0.00

Target Compounds Qvalue

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

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