

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010824.D
 Acq On : 25 Jun 2022 12:47
 Operator : CG/JU
 Sample : PB145662BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK662

Quant Time: Jun 26 22:51:48 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	7.722	152	433785	20.000	ng/u1	0.00
20) Naphthalene-d8	10.516	136	1738058	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.375	164	923384	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.139	188	1698814	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.251	240	1661364	20.000	ng/u1	# 0.00
88) Perylene-d12	23.621	264	1763452	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	63619	5.548	ng/uL	0.00
4) Pyridine-d5	3.617	84	824676	27.017	ng/u1	0.01
7) Phenol-d5	6.905	99	956414	27.852	ng/u1	0.01
9) Bis-(2-Chloroethyl)eth...	7.064	67	617000	27.474	ng/u1	0.00
11) 2-Chlorophenol-d4	7.258	132	827661	29.051	ng/u1	0.00
15) 4-Methylphenol-d8	8.440	113	759871	27.993	ng/u1	0.00
21) Nitrobenzene-d5	8.881	128	371512	33.577	ng/u1	0.00
24) 2-Nitrophenol-d4	9.599	143	360168	39.346	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.140	165	633191	27.995	ng/u1	0.00
31) 4-Chloroaniline-d4	10.663	131	1031926	27.950	ng/u1	0.00
46) Dimethylphthalate-d6	13.798	166	1830933	28.947	ng/u1	0.00
49) Acenaphthylene-d8	14.069	160	2222840	27.535	ng/u1	0.00
54) 4-Nitrophenol-d4	14.598	143	272686	27.601	ng/u1	0.02
60) Fluorene-d10	15.375	176	1539546	28.363	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.493	200	127421	22.639	ng/u1	0.00
73) Anthracene-d10	17.239	188	2262277	29.931	ng/u1	0.00
81) Pyrene-d10	19.486	212	2515722	27.794	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.469	264	2667061	30.606	ng/u1	0.00

Target Compounds Qvalue

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

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