

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
 Data File : BP009708.D
 Acq On : 08 Apr 2022 09:41
 Operator : CG/JU
 Sample : PB143863BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK863

Quant Time: Apr 08 22:25:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 07 19:16:52 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	115442	20.000	ng/ul	0.00
20) Naphthalene-d8	10.786	136	512724	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.610	164	378940	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.363	188	866723	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.445	240	951225	20.000	ng/ul	# 0.00
88) Perylene-d12	23.921	264	923658	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	15744	6.087	ng/uL	0.00
4) Pyridine-d5	3.810	84	210652	29.537	ng/ul	0.00
7) Phenol-d5	7.128	99	268265	26.469	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.298	67	178560	29.593	ng/ul	0.00
11) 2-Chlorophenol-d4	7.504	132	218367	28.563	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	230484	27.242	ng/ul	0.00
21) Nitrobenzene-d5	9.134	128	112626	30.461	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	121847	30.059	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	224386	26.453	ng/ul	0.00
31) 4-Chloroaniline-d4	10.916	131	340312	28.168	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	927659	31.165	ng/ul	0.00
49) Acenaphthylene-d8	14.304	160	974787	27.642	ng/ul	0.00
54) 4-Nitrophenol-d4	14.780	143	149427	27.680	ng/ul	0.00
60) Fluorene-d10	15.604	176	749251	29.922	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.710	200	130132	25.345	ng/ul	0.00
73) Anthracene-d10	17.463	188	1253960	30.217	ng/ul	0.00
81) Pyrene-d10	19.686	212	1577798	30.371	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.762	264	1514087	31.477	ng/ul	0.00

Target Compounds Qvalue

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

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