

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030322\
 Data File : BP009252.D
 Acq On : 03 Mar 2022 09:50
 Operator : CG/JU
 Sample : PB142783BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK783

Quant Time: Mar 03 11:03:52 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030222.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 02 18:15:01 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	347312	20.000	ng/u1	# 0.00
20) Naphthalene-d8	10.616	136	1400307	20.000	ng/u1	#-0.01
38) Acenaphthene-d10	14.463	164	871228	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.222	188	1797380	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.327	240	1808696	20.000	ng/u1	-0.01
88) Perylene-d12	23.739	264	1850700	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.323	96	51835	5.999	ng/uL	0.00
4) Pyridine-d5	3.723	84	668994	31.028	ng/u1	0.00
7) Phenol-d5	6.987	99	880072	31.280	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67	534379	31.648	ng/u1	0.00
11) 2-Chlorophenol-d4	7.358	132	704554	32.030	ng/u1	0.00
15) 4-Methylphenol-d8	8.528	113	673053	31.295	ng/u1	0.00
21) Nitrobenzene-d5	8.975	128	287312	35.182	ng/u1	0.00
24) 2-Nitrophenol-d4	9.699	143	289056	37.345	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	642732	31.165	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.752	131	933142	31.998	ng/u1	0.00
46) Dimethylphthalate-d6	13.875	166	1997060	32.708	ng/u1	-0.01
49) Acenaphthylene-d8	14.151	160	2501021	31.001	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.645	143	251138	25.972	ng/u1	-0.01
60) Fluorene-d10	15.457	176	1728866	32.542	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.569	200	169079	23.176	ng/u1	-0.01
73) Anthracene-d10	17.322	188	2691389	31.861	ng/u1	-0.01
81) Pyrene-d10	19.569	212	3124351	31.425	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.580	264	3037295	32.448	ng/u1	-0.02

Target Compounds Qvalue

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

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