

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006919.D
 Acq On : 10 Sep 2021 02:39
 Operator : CG/JU
 Sample : PB138845BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK845

Quant Time: Sep 10 04:18:07 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 09 11:07:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	379886	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	1527975	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.563	164	944806	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.310	188	1945779	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.392	240	1859959	20.000	ng/u1	# 0.00
88) Perylene-d12	23.827	264	1931092	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	59483	6.231	ng/uL	0.00
4) Pyridine-d5	3.787	84	783233	31.485	ng/u1	0.00
7) Phenol-d5	7.093	99	934030	33.453	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	542711	33.844	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	753500	33.429	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	714497	32.587	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	334636	36.157	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	340138	37.330	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.351	165	675612	31.446	ng/u1	0.00
31) 4-Chloroaniline-d4	10.869	131	987355	33.231	ng/u1	0.00
46) Dimethylphthalate-d6	13.969	166	2150281	35.139	ng/u1	0.00
49) Acenaphthylene-d8	14.251	160	2573978	33.212	ng/u1	0.00
54) 4-Nitrophenol-d4	14.739	143	322804	30.364	ng/u1	0.00
60) Fluorene-d10	15.551	176	1828371	33.854	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	242005	28.115	ng/u1	0.00
73) Anthracene-d10	17.410	188	2963360	36.786	ng/u1	0.00
81) Pyrene-d10	19.639	212	3399265	39.024	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.668	264	3282354	35.233	ng/u1	-0.01

Target Compounds	Qvalue

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

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