

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006904.D
 Acq On : 09 Sep 2021 17:33
 Operator : CG/JU
 Sample : PB138738BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK738

Quant Time: Sep 10 01:04:31 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	7.934	152	280636	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	1119283	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.563	164	685074	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.310	188	1419729	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.392	240	1380516	20.000	ng/u1	0.00
88) Perylene-d12	23.833	264	1441452	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	45571	6.462	ng/uL	0.00
4) Pyridine-d5	3.781	84	586924	31.938	ng/u1	0.00
7) Phenol-d5	7.087	99	685245	33.222	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	410768	34.676	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	553260	33.226	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	519286	32.060	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	244298	36.034	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	237220	35.541	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	484994	30.816	ng/u1	0.00
31) 4-Chloroaniline-d4	10.869	131	719437	33.055	ng/u1	0.00
46) Dimethylphthalate-d6	13.975	166	1564217	35.253	ng/u1	0.00
49) Acenaphthylene-d8	14.257	160	1865362	33.194	ng/u1	0.00
54) 4-Nitrophenol-d4	14.745	143	221885	28.785	ng/u1	0.00
60) Fluorene-d10	15.557	176	1333354	34.049	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	168225	26.785	ng/u1	0.00
73) Anthracene-d10	17.410	188	2164420	36.824	ng/u1	0.00
81) Pyrene-d10	19.645	212	2515366	38.906	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.674	264	2440412	35.094	ng/u1	0.00

Target Compounds	Qvalue

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

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