

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006907.D
 Acq On : 09 Sep 2021 19:16
 Operator : CG/JU
 Sample : PB138901BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK901

Quant Time: Sep 10 01:05:17 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	308380	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	1241892	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.563	164	754261	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.310	188	1536063	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.392	240	1461322	20.000	ng/ul	# 0.00
88) Perylene-d12	23.827	264	1508105	20.000	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	50450	6.510	ng/uL	0.00
4) Pyridine-d5	3.787	84	654109	32.392	ng/ul	0.00
7) Phenol-d5	7.093	99	764545	33.732	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	454326	34.902	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	615864	33.658	ng/ul	0.00
15) 4-Methylphenol-d8	8.634	113	582533	32.729	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	267900	35.614	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	258009	34.839	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	538367	30.830	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	787989	32.631	ng/ul	0.00
46) Dimethylphthalate-d6	13.969	166	1705374	34.909	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	2053774	33.194	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	243870	28.735	ng/ul	0.00
60) Fluorene-d10	15.551	176	1460142	33.866	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	179058	26.351	ng/ul	0.00
73) Anthracene-d10	17.410	188	2340993	36.812	ng/ul	0.00
81) Pyrene-d10	19.639	212	2691413	39.327	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.669	264	2582641	35.498	ng/ul	-0.01

Target Compounds	Qvalue

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

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