

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006920.D
 Acq On : 10 Sep 2021 03:13
 Operator : CG/JU
 Sample : PB138946BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK946

Quant Time: Sep 10 04:18:22 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	372323	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	1507555	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.563	164	930117	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.310	188	1878124	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.386	240	1771908	20.000	ng/u1	-0.01
88) Perylene-d12	23.821	264	1812242	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	59247	6.333	ng/uL	0.01
4) Pyridine-d5	3.787	84	766179	31.425	ng/u1	0.00
7) Phenol-d5	7.093	99	920118	33.624	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	546944	34.801	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	744687	33.709	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	710530	33.064	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	334290	36.609	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	330995	36.818	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	669565	31.586	ng/u1	0.00
31) 4-Chloroaniline-d4	10.869	131	977746	33.354	ng/u1	0.00
46) Dimethylphthalate-d6	13.969	166	2094702	34.772	ng/u1	0.00
49) Acenaphthylene-d8	14.251	160	2535265	33.229	ng/u1	0.00
54) 4-Nitrophenol-d4	14.740	143	311327	29.747	ng/u1	0.00
60) Fluorene-d10	15.551	176	1783493	33.545	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	230470	27.740	ng/u1	0.00
73) Anthracene-d10	17.404	188	2862746	36.818	ng/u1	-0.01
81) Pyrene-d10	19.639	212	3272755	39.439	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.663	264	3103118	35.494	ng/u1	-0.02

Target Compounds	Qvalue

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

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