

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006930.D
 Acq On : 10 Sep 2021 08:55
 Operator : CG/JU
 Sample : PB138946BL
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK946

Quant Time: Sep 10 10:45:05 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	328257	20.000	ng/ul	0.00
20) Naphthalene-d8	10.734	136	1314981	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.557	164	808902	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.310	188	1650414	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.392	240	1586865	20.000	ng/ul	# 0.00
88) Perylene-d12	23.827	264	1624905	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	48010	5.820	ng/uL	0.00
4) Pyridine-d5	3.787	84	623682	29.015	ng/ul	0.00
7) Phenol-d5	7.093	99	720899	29.880	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	427979	30.887	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	585188	30.045	ng/ul	0.00
15) 4-Methylphenol-d8	8.634	113	555714	29.332	ng/ul	0.00
21) Nitrobenzene-d5	9.093	128	262278	32.929	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	258891	33.015	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.351	165	527011	28.502	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	772900	30.227	ng/ul	0.00
46) Dimethylphthalate-d6	13.969	166	1683963	32.142	ng/ul	0.00
49) Acenaphthylene-d8	14.251	160	2031032	30.609	ng/ul	0.00
54) 4-Nitrophenol-d4	14.739	143	236700	26.006	ng/ul	0.00
60) Fluorene-d10	15.551	176	1452162	31.406	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	177269	24.280	ng/ul	0.00
73) Anthracene-d10	17.410	188	2292847	33.557	ng/ul	0.00
81) Pyrene-d10	19.639	212	2605230	35.056	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264	2457800	31.354	ng/ul	-0.01

Target Compounds Qvalue

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

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