

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006905.D
 Acq On : 09 Sep 2021 18:07
 Operator : CG/JU
 Sample : PB138916BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK916

Quant Time: Sep 10 01:04:46 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	307285	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	1222848	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.563	164	746940	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.310	188	1529278	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.392	240	1441813	20.000	ng/u1	# 0.00
88) Perylene-d12	23.827	264	1494576	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	49096	6.358	ng/uL	0.00
4) Pyridine-d5	3.787	84	633655	31.490	ng/u1	0.00
7) Phenol-d5	7.093	99	745545	33.011	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	444695	34.284	ng/u1	0.00
11) 2-Chlorophenol-d4	7.469	132	604876	33.176	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	573455	32.334	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	266965	36.043	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	256681	35.199	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	535608	31.150	ng/u1	0.00
31) 4-Chloroaniline-d4	10.869	131	779668	32.789	ng/u1	0.00
46) Dimethylphthalate-d6	13.969	166	1672553	34.573	ng/u1	0.00
49) Acenaphthylene-d8	14.257	160	2016996	32.919	ng/u1	0.00
54) 4-Nitrophenol-d4	14.740	143	241460	28.730	ng/u1	0.00
60) Fluorene-d10	15.551	176	1447043	33.891	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	177492	26.236	ng/u1	0.00
73) Anthracene-d10	17.410	188	2308588	36.463	ng/u1	0.00
81) Pyrene-d10	19.639	212	2640659	39.107	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.669	264	2539230	35.217	ng/u1	-0.01

Target Compounds Qvalue

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

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