

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006906.D
 Acq On : 09 Sep 2021 18:42
 Operator : CG/JU
 Sample : PB138900BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK900

Quant Time: Sep 10 01:05:02 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	313787	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	1254608	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.563	164	776792	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.310	188	1609385	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	1532470	20.000	ng/ul	-0.01
88) Perylene-d12	23.827	264	1596250	20.000	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	51855	6.576	ng/uL	0.00
4) Pyridine-d5	3.787	84	652280	31.744	ng/ul	0.00
7) Phenol-d5	7.093	99	765623	33.198	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	456335	34.453	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	618139	33.201	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	588812	32.512	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	272773	35.894	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	263338	35.198	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	544976	30.892	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	812107	33.289	ng/ul	0.00
46) Dimethylphthalate-d6	13.969	166	1774424	35.269	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	2111670	33.140	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	256210	29.313	ng/ul	0.00
60) Fluorene-d10	15.551	176	1518662	34.202	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	189833	26.664	ng/ul	0.00
73) Anthracene-d10	17.410	188	2437344	36.581	ng/ul	0.00
81) Pyrene-d10	19.639	212	2817301	39.255	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.669	264	2690355	34.936	ng/ul	-0.01

Target Compounds	Qvalue

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

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