

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006903.D
 Acq On : 09 Sep 2021 16:47
 Operator : CG/JU
 Sample : PB138799BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK799

Quant Time: Sep 10 01:03:25 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	283688	20.000	ng/ul	0.00
20) Naphthalene-d8	10.739	136	1147288	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.563	164	728398	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.310	188	1548102	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.392	240	1520948	20.000	ng/ul	0.00
88) Perylene-d12	23.827	264	1545233	20.000	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	45334	6.359	ng/uL	0.00
4) Pyridine-d5	3.781	84	581107	31.281	ng/ul	0.00
7) Phenol-d5	7.087	99	685437	32.874	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	407260	34.010	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	555552	33.005	ng/ul	0.00
15) 4-Methylphenol-d8	8.634	113	534158	32.623	ng/ul	0.00
21) Nitrobenzene-d5	9.092	128	253236	36.441	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	247260	36.141	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.351	165	501643	31.096	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	749802	33.610	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	1690414	35.831	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1985551	33.231	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	250877	30.610	ng/ul	0.00
60) Fluorene-d10	15.557	176	1431539	34.382	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	191934	28.026	ng/ul	0.00
73) Anthracene-d10	17.410	188	2351677	36.692	ng/ul	0.00
81) Pyrene-d10	19.639	212	2742280	38.499	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.674	264	2633746	35.330	ng/ul	0.00

Target Compounds	Qvalue

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

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