

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006922.D
 Acq On : 10 Sep 2021 04:21
 Operator : CG/JU
 Sample : PB138976BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK976

Quant Time: Sep 10 06:32:51 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	348280	20.000	ng/ul	0.00
20) Naphthalene-d8	10.734	136	1402779	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.557	164	876152	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.310	188	1824661	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	1798170	20.000	ng/ul	-0.01
88) Perylene-d12	23.822	264	1867895	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	56130	6.414	ng/uL	0.00
4) Pyridine-d5	3.781	84	719228	31.536	ng/ul	0.00
7) Phenol-d5	7.093	99	857163	33.486	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.264	67	513633	34.938	ng/ul	0.00
11) 2-Chlorophenol-d4	7.464	132	692190	33.496	ng/ul	0.00
15) 4-Methylphenol-d8	8.634	113	660879	32.877	ng/ul	0.00
21) Nitrobenzene-d5	9.093	128	304591	35.848	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	301024	35.985	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	614470	31.152	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	914642	33.531	ng/ul	0.00
46) Dimethylphthalate-d6	13.969	166	1991819	35.100	ng/ul	0.00
49) Acenaphthylene-d8	14.251	160	2383512	33.164	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	300991	30.531	ng/ul	0.00
60) Fluorene-d10	15.551	176	1706620	34.076	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	220733	27.346	ng/ul	0.00
73) Anthracene-d10	17.404	188	2776917	36.760	ng/ul	-0.01
81) Pyrene-d10	19.639	212	3233784	38.400	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.669	264	3158796	35.054	ng/ul	-0.01

Target Compounds Qvalue

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

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