

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021822\
 Data File : BP009227.D
 Acq On : 18 Feb 2022 22:37
 Operator : CG/JU
 Sample : N1466-10DL 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A5-10-4.0DL

Quant Time: Feb 19 00:51:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 19 00:46:08 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	257117	20.000	ng	0.00
21) Naphthalene-d8	10.557	136	1038280	20.000	ng	0.00
39) Acenaphthene-d10	14.410	164	734536	20.000	ng	0.00
64) Phenanthrene-d10	17.169	188	1585017	20.000	ng	0.00
76) Chrysene-d12	21.275	240	1716631	20.000	ng	0.00
86) Perylene-d12	23.657	264	1522099	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	88774	6.159	ng	0.01
7) Phenol-d6	6.981	99	123925	6.525	ng	0.02
23) Nitrobenzene-d5	8.940	82	90360	4.908	ng	0.00
42) 2,4,6-Tribromophenol	15.928	330	50684	5.168	ng	0.00
45) 2-Fluorobiphenyl	13.028	172	283393	5.401	ng	0.00
79) Terphenyl-d14	19.739	244	488552	5.118	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.610	128	2404845	45.412	ng	98
37) 2-Methylnaphthalene	12.228	142	2512489	67.046	ng	98
38) 1-Methylnaphthalene	12.446	142	1274000	34.792	ng	99
46) 1,1'-Biphenyl	13.240	154	173977	3.224	ng	97
58) Fluorene	15.463	166	137866	2.708	ng	# 87
71) Phenanthrene	17.216	178	409727	4.774	ng	97

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

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