

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121021\
 Data File : BP008362.D
 Acq On : 11 Dec 2021 03:19
 Operator : CG/JU
 Sample : M4964-02 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EAST-SIDEWALL

Quant Time: Dec 11 04:04:01 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 11 03:53:01 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.758	152	65769	20.000	ng	0.00
21) Naphthalene-d8	10.557	136	234082	20.000	ng	0.00
39) Acenaphthene-d10	14.416	164	153408	20.000	ng	0.00
64) Phenanthrene-d10	17.181	188	347135	20.000	ng	0.00
76) Chrysene-d12	21.292	240	401340	20.000	ng	0.00
86) Perylene-d12	23.680	264	437964	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.364	112	12316	3.015	ng	0.00
7) Phenol-d6	6.969	99	13983	2.536	ng	0.02
23) Nitrobenzene-d5	8.934	82	10161	1.974	ng	0.00
42) 2,4,6-Tribromophenol	15.922	330	6280	3.203	ng	0.00
45) 2-Fluorobiphenyl	13.028	172	22335	2.115	ng	0.00
79) Terphenyl-d14	19.751	244	44604	2.323	ng	0.00
Target Compounds						
					Qvalue	
35) Caprolactam	11.593	113	4663	5.456	ng	# 1
37) 2-Methylnaphthalene	12.222	142	31698	3.737	ng	# 96
38) 1-Methylnaphthalene	12.446	142	27504	3.332	ng	# 95
71) Phenanthrene	17.222	178	103693	5.658	ng	# 94
83) Chrysene	21.327	228	79361	3.135	ng	# 52

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

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