

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010322\
 Data File : BP008646.D
 Acq On : 03 Jan 2022 19:56
 Operator : CG/JU
 Sample : M5257-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P2-D

Quant Time: Jan 04 06:01:20 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP122821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 30 08:44:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	582776	20.000	ng	-0.01
21) Naphthalene-d8	10.687	136	2199481	20.000	ng	-0.01
39) Acenaphthene-d10	14.516	164	1359696	20.000	ng	-0.01
64) Phenanthrene-d10	17.269	188	2621791	20.000	ng	0.00
76) Chrysene-d12	21.351	240	2277793	20.000	ng	-0.01
86) Perylene-d12	23.774	264	2442424	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.470	112	4182647	131.001	ng	0.00
7) Phenol-d6	7.058	99	4685649	105.435	ng	0.00
23) Nitrobenzene-d5	9.046	82	3410496	92.243	ng	-0.01
42) 2,4,6-Tribromophenol	16.010	330	2559427	139.128	ng	0.00
45) 2-Fluorobiphenyl	13.146	172	8832700	94.116	ng	-0.01
79) Terphenyl-d14	19.828	244	10614152	95.144	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.740	128	4282817	37.515	ng	100
32) Benzoic acid	9.952	122	142129	7.456	ng	96
37) 2-Methylnaphthalene	12.346	142	176767	2.057	ng	99

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

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