

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121321\
 Data File : BP008388.D
 Acq On : 13 Dec 2021 21:48
 Operator : CG/JU
 Sample : M5051-01DL2 25X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P-1DL2

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 12/14/2021
 Supervised By :Jagrut Upadhyay 12/14/2021

Quant Time: Dec 14 02:21:37 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	7.752	152	75851	20.000	ng	-0.01
21) Naphthalene-d8	10.546	136	280497	20.000	ng	-0.01
39) Acenaphthene-d10	14.404	164	187354	20.000	ng	-0.01
64) Phenanthrene-d10	17.169	188	399976	20.000	ng	#-0.01
76) Chrysene-d12	21.274	240	451943	20.000	ng	-0.01
86) Perylene-d12	23.651	264	503776	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.358	112	14206	3.016	ng	0.00
7) Phenol-d6	6.963	99	17085	2.687	ng	0.01
23) Nitrobenzene-d5	8.916	82	14506	2.352	ng	-0.01
42) 2,4,6-Tribromophenol	15.916	330	7520	3.140	ng	0.00
45) 2-Fluorobiphenyl	13.022	172	29726	2.305	ng	-0.01
79) Terphenyl-d14	19.739	244	52445	2.425	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	10.593	128	838288	58.340	ng	99
37) 2-Methylnaphthalene	12.216	142	79491	7.822	ng	98
38) 1-Methylnaphthalene	12.434	142	41316m	4.177	ng	
71) Phenanthrene	17.210	178	136707	6.474	ng	99
75) Fluoranthene	19.192	202	64530	2.268	ng	98

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

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