

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012720\
 Data File : BG044220.D
 Acq On : 27 Jan 2020 18:09
 Operator : CG/JU
 Sample : L1245-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20200065-AMENDED-COMP-3

Manual Integrations
 APPROVED

mohammad
 1/28/2020 1:12:39 PM

Quant Time: Jan 28 01:17:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 31 13:32:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	85532	20.00	ng	-0.04
21) Naphthalene-d8	10.73	136	386388	20.00	ng	-0.04
39) Acenaphthene-d10	14.56	164	258338	20.00	ng	-0.04
64) Phenanthrene-d10	17.31	188	502534	20.00	ng	-0.03
76) Chrysene-d12	21.55	240	444982	20.00	ng	-0.04
87) Perylene-d12	24.66	264	499866	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	300404	64.49	ng	-0.03
7) Phenol-d6	7.10	99	247900	38.07	ng	-0.02
23) Nitrobenzene-d5	9.09	82	592005	81.96	ng	-0.04
42) 2,4,6-Tribromophenol	16.06	330	382492	141.48	ng	-0.03
45) 2-Fluorobiphenyl	13.19	172	1389354	97.44	ng	-0.04
79) Terphenyl-d14	19.93	244	1783615	101.55	ng	-0.03
Target Compounds						
10) Phenol	7.12	94	19758	2.945	ng	92
20) 3+4-Methylphenols	8.70	107	90138	13.309	ng	92
32) Benzoic acid	10.05	122	3997m	7.100	ng	
50) Dimethylphthalate	14.02	163	94967	5.177	ng	97
52) Acenaphthene	14.63	154	31948	2.118	ng	97

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

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