

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG040720\
 Data File : BG044963.D
 Acq On : 7 Apr 2020 20:20
 Operator : CG/JU
 Sample : L2161-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RECLAIM-SYSTEM-SEDIMENT-PIT

Manual Integrations
 APPROVED

mohammad
 4/8/2020 8:14:17 AM

Quant Time: Apr 08 07:18:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG031020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 23 07:18:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	86611	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	326619	20.00	ng	0.00
39) Acenaphthene-d10	14.68	164	217784	20.00	ng	0.00
64) Phenanthrene-d10	17.42	188	503756	20.00	ng	-0.01
76) Chrysene-d12	21.68	240	511108	20.00	ng	-0.02
87) Perylene-d12	24.90	264	554627	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	549252	98.72	ng	0.00
7) Phenol-d6	7.20	99	728817	90.16	ng	0.00
23) Nitrobenzene-d5	9.19	82	585109	78.61	ng	0.00
42) 2,4,6-Tribromophenol	16.16	330	493184	172.95	ng	-0.01
45) 2-Fluorobiphenyl	13.30	172	1331552	87.56	ng	0.00
79) Terphenyl-d14	20.03	244	2180374	79.80	ng	0.00
Target Compounds						
10) Phenol	7.23	94	59959	7.492	ng	91
20) 3+4-Methylphenols	8.80	107	367986	49.241	ng	90
32) Benzoic acid	10.10	122	8188m	11.444	ng	
50) Dimethylphthalate	14.12	163	64331	3.709	ng	# 98

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

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