

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111519\
 Data File : BG043670.D
 Acq On : 16 Nov 2019 3:39
 Operator : JU
 Sample : K5837-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EFF-WW

Manual Integrations
 APPROVED

mohammad

11/18/2019 3:07:20 PM

Quant Time: Nov 18 07:06:27 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 04 15:22:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	28789	20.00	ng	-0.01
21) Naphthalene-d8	10.80	136	108310	20.00	ng	-0.02
39) Acenaphthene-d10	14.62	164	79004	20.00	ng	-0.02
64) Phenanthrene-d10	17.37	188	180267	20.00	ng	-0.02
76) Chrysene-d12	21.64	240	206677	20.00	ng	-0.02
87) Perylene-d12	24.81	264	238996	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	89562	57.01	ng	0.00
7) Phenol-d6	7.20	99	93180	41.22	ng	0.00
23) Nitrobenzene-d5	9.16	82	181736	86.50	ng	-0.02
42) 2,4,6-Tribromophenol	16.12	330	132581	88.18	ng	-0.01
45) 2-Fluorobiphenyl	13.25	172	521672	91.24	ng	-0.02
79) Terphenyl-d14	19.99	244	958041	86.96	ng	-0.01
Target Compounds						
10) Phenol	7.23	94	31322	15.164	ng	95
20) 3+4-Methylphenols	8.79	107	4787	2.318	ng	93
32) Benzoic acid	10.50	122	50341m	47.682	ng	
50) Dimethylphthalate	14.07	163	74164	12.119	ng	99

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

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