

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062220\
 Data File : BG045827.D
 Acq On : 22 Jun 2020 14:36
 Operator : CG/JU
 Sample : L3033-06DL 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 KY032MW006-200616DL

Manual Integrations
 APPROVED

Sohil
 6/23/2020 2:19:10 PM

Quant Time: Jun 22 15:25:26 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 15 18:34:19 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	52222	20.00	ng	-0.01
21) Naphthalene-d8	10.72	136	193537	20.00	ng	0.00
39) Acenaphthene-d10	14.57	164	142931	20.00	ng	0.00
64) Phenanthrene-d10	17.32	188	326528	20.00	ng	0.00
76) Chrysene-d12	21.57	240	357760	20.00	ng	0.00
86) Perylene-d12	24.70	264	375958	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	17109	5.53	ng	-0.01
7) Phenol-d6	7.15	99	19926	4.24	ng	-0.02
23) Nitrobenzene-d5	9.08	82	36797	8.69	ng	0.00
42) 2,4,6-Tribromophenol	16.06	330	27358	12.68	ng	0.00
45) 2-Fluorobiphenyl	13.19	172	89660	9.22	ng	0.00
79) Terphenyl-d14	19.93	244	165814	9.39	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.36	88	709633	501.186	ng	96
10) Phenol	7.17	94	82400	18.085	ng	99
11) bis(2-Chloroethyl)ether	7.33	93	134545	38.877	ng	96
13) 1,4-Dichlorobenzene	7.95	146	26965	6.846	ng	93
14) 1,2-Dichlorobenzene	8.27	146	127537	33.822	ng	98
17) 2-Methylphenol	8.40	107	116336	34.309	ng	97
20) 3+4-Methylphenols	8.74	107	764578	171.451	ng	# 73
27) 2,4-Dimethylphenol	9.97	122	313064m	108.961	ng	
31) Naphthalene	10.77	128	50821	4.803	ng	98

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

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