

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062220\
 Data File : BG045826.D
 Acq On : 22 Jun 2020 13:56
 Operator : CG/JU
 Sample : L3033-10DL2 50X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 Z1-DUP-0572-200616DL2

Manual Integrations
 APPROVED

Sohil
 6/23/2020 2:19:09 PM

Quant Time: Jun 22 14:33:13 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	42568	20.00	ng	0.00
21) Naphthalene-d8	10.72	136	169962	20.00	ng	0.00
39) Acenaphthene-d10	14.56	164	132155	20.00	ng	0.00
64) Phenanthrene-d10	17.31	188	320580	20.00	ng	0.00
76) Chrysene-d12	21.57	240	335967	20.00	ng	0.00
86) Perylene-d12	24.70	264	350073	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	2575	1.02	ng	0.00
7) Phenol-d6	7.15	99	2380	0.62	ng	-0.01
23) Nitrobenzene-d5	9.08	82	5388	1.45	ng	0.00
42) 2,4,6-Tribromophenol	16.07	330	4234	2.12	ng	0.00
45) 2-Fluorobiphenyl	13.18	172	14762	1.64	ng	0.00
79) Terphenyl-d14	19.93	244	28652	1.73	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.36	88	123846	107.304	ng	99
10) Phenol	7.17	94	14308	3.852	ng	98
11) bis(2-Chloroethyl)ether	7.34	93	21841	7.742	ng	95
14) 1,2-Dichlorobenzene	8.27	146	21589	7.024	ng	89
17) 2-Methylphenol	8.41	107	20977	7.589	ng	93
20) 3+4-Methylphenols	8.74	107	141331	38.880	ng	# 74
27) 2,4-Dimethylphenol	9.96	122	57610m	22.832	ng	

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

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