

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG060520\
 Data File : BG045615.D
 Acq On : 5 Jun 2020 13:56
 Operator : CG/JU
 Sample : L2822-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AP7

Manual Integrations
 APPROVED

mohammad
 6/8/2020 12:48:52 PM

Quant Time: Jun 06 05:27:35 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG051520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 01 15:07:00 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.95	152	68926	20.00	ng	0.00
21) Naphthalene-d8	10.77	136	306080	20.00	ng	0.01
39) Acenaphthene-d10	14.61	164	191123	20.00	ng	0.02
64) Phenanthrene-d10	17.36	188	371251	20.00	ng	0.02
76) Chrysene-d12	21.61	240	372773	20.00	ng	0.00
87) Perylene-d12	24.76	264	408956	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	530792	150.50	ng	0.01
7) Phenol-d6	7.19	99	657531	130.99	ng	0.05
23) Nitrobenzene-d5	9.14	82	899553	160.27	ng	0.03
42) 2,4,6-Tribromophenol	16.12	330	407002	166.51	ng	0.03
45) 2-Fluorobiphenyl	13.24	172	1249605	110.90	ng	0.03
79) Terphenyl-d14	19.96	244	1931139	120.82	ng	0.00
Target Compounds						
3) Pyridine	3.85	79	13183m	2.71	ng	Qvalue
10) Phenol	7.22	94	1945047	375.96	ng	93
17) 2-Methylphenol	8.47	107	2310246	596.59	ng	94
20) 3+4-Methylphenols	8.83	107	4648741	881.58	ng	85
27) 2,4-Dimethylphenol	10.02	122	3121677	818.16	ng	99
31) Naphthalene	10.82	128	139749	9.80	ng	98
37) 2-Methylnaphthalene	12.43	142	64247	6.04	ng	# 90
38) 1-Methylnaphthalene	12.65	142	42073	4.19	ng	# 98

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

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