

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061120\
 Data File : BG045735.D
 Acq On : 12 Jun 2020 00:21
 Operator : CG/JU
 Sample : L2973-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RL-356-RL-212

Manual Integrations
 APPROVED

mohammad
 6/15/2020 8:44:55 AM

Quant Time: Jun 12 03:41:45 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	70441	20.00	ng	0.00
21) Naphthalene-d8	10.75	136	282952	20.00	ng	-0.01
39) Acenaphthene-d10	14.58	164	189069	20.00	ng	-0.01
64) Phenanthrene-d10	17.33	188	396626	20.00	ng	-0.01
76) Chrysene-d12	21.59	240	354758	20.00	ng	-0.01
87) Perylene-d12	24.74	264	402179	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	19522	5.42	ng	0.00
7) Phenol-d6	7.13	99	140628	27.41	ng	-0.02
23) Nitrobenzene-d5	9.10	82	158783	30.60	ng	-0.01
42) 2,4,6-Tribromophenol	16.10	330	1014	0.42	ng	0.00
45) 2-Fluorobiphenyl	13.21	172	373498	33.51	ng	0.00
79) Terphenyl-d14	19.95	244	534806	35.16	ng	-0.01
Target Compounds						
50) Dimethylphthalate	14.04	163	62917	4.851	ng	Qvalue # 99
75) Fluoranthene	19.39	202	188541	8.561	ng	98
78) Pyrene	19.75	202	347342	17.176	ng	98
81) Benzo(a)anthracene	21.57	228	229557m	11.616	ng	
83) Chrysene	21.64	228	225861m	11.886	ng	
86) Indeno(1,2,3-cd)pyrene	28.30	276	330822	13.313	ng	97
88) Benzo(b)fluoranthene	23.74	252	442394m	20.510	ng	
89) Benzo(k)fluoranthene	23.80	252	113634m	5.608	ng	
90) Benzo(a)pyrene	24.58	252	457030	22.751	ng	98
91) Dibenzo(a,h)anthracene	28.33	278	61293	3.036	ng	# 94
92) Benzo(g,h,i)perylene	29.41	276	432248	21.410	ng	97

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

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