

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG042020\
 Data File : BG045093.D
 Acq On : 20 Apr 2020 13:06
 Operator : CG/JU
 Sample : L2323-12
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-19-COMP

Manual Integrations
 APPROVED

mohammad
 4/21/2020 2:53:53 PM

Quant Time: Apr 20 14:00:46 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 20 11:57:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	72072	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	317043	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	234648	20.00	ng	0.00
64) Phenanthrene-d10	17.40	188	516153	20.00	ng	0.00
76) Chrysene-d12	21.67	240	465175	20.00	ng	0.00
87) Perylene-d12	24.86	264	510011	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	466311	106.82	ng	0.00
7) Phenol-d6	7.19	99	769073	118.57	ng	0.00
23) Nitrobenzene-d5	9.18	82	568710	81.85	ng	0.00
42) 2,4,6-Tribromophenol	16.15	330	445939	123.02	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	1416344	88.51	ng	0.00
79) Terphenyl-d14	20.02	244	2051195	96.11	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.11	163	339612	17.775	ng	99
71) Phenanthrene	17.45	178	61589	2.322	ng	98
78) Pyrene	19.81	202	80245	2.502	ng	96
81) Benzo(a)anthracene	21.65	228	62710	2.080	ng	96
83) Chrysene	21.71	228	66442	2.294	ng	98
86) Indeno(1,2,3-cd)pyrene	28.48	276	77823	2.023	ng	# 86
88) Benzo(b)fluoranthene	23.85	252	94523m	2.959	ng	
90) Benzo(a)pyrene	24.71	252	77862	2.581	ng	100
92) Benzo(g,h,i)perylene	29.60	276	86512	2.839	ng	# 92

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

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