

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG083120\
 Data File : BG046480.D
 Acq On : 31 Aug 2020 22:33
 Operator : CG/JU
 Sample : L3847-25
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LINDEN-JOP-BH-12-A

Manual Integrations
 APPROVED

mohammad
 9/1/2020 10:47:50 AM

Quant Time: Sep 01 04:51:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 31 16:05:33 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	88253	20.00	ng	0.00
21) Naphthalene-d8	10.73	136	379962	20.00	ng	0.00
39) Acenaphthene-d10	14.56	164	251128	20.00	ng	0.00
64) Phenanthrene-d10	17.31	188	533471	20.00	ng	0.00
76) Chrysene-d12	21.55	240	496775	20.00	ng	0.00
86) Perylene-d12	24.66	264	520228	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	518900	104.14	ng	0.00
7) Phenol-d6	7.11	99	718011	101.65	ng	0.00
23) Nitrobenzene-d5	9.09	82	483343	72.71	ng	0.00
42) 2,4,6-Tribromophenol	16.05	330	306401	90.05	ng	0.00
45) 2-Fluorobiphenyl	13.19	172	1245768	75.73	ng	0.00
79) Terphenyl-d14	19.92	244	1517299	65.00	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.01	163	294654	15.166	ng	99
71) Phenanthrene	17.35	178	179115	6.624	ng	98
75) Fluoranthene	19.36	202	463310	13.314	ng	97
78) Pyrene	19.72	202	428530	13.546	ng	98
81) Benzo(a)anthracene	21.54	228	257627	8.320	ng	98
83) Chrysene	21.60	228	233079	7.825	ng	99
87) Indeno(1,2,3-cd)pyrene	28.16	276	175560	4.699	ng	97
88) Benzo(b)fluoranthene	23.68	252	337456	10.388	ng	# 98
89) Benzo(k)fluoranthene	23.73	252	103805m	3.307	ng	
90) Benzo(a)pyrene	24.51	252	246664	8.137	ng	98
92) Benzo(g,h,i)perylene	29.25	276	178500	5.928	ng	99

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

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