

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG083120\
 Data File : BG046479.D
 Acq On : 31 Aug 2020 21:53
 Operator : CG/JU
 Sample : L3847-19
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LINDEN-JOP-BH-11-A

Manual Integrations
 APPROVED

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

Quant Time: Sep 01 04:41:10 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	90688	20.00	ng	0.00
21) Naphthalene-d8	10.73	136	385712	20.00	ng	0.00
39) Acenaphthene-d10	14.56	164	263282	20.00	ng	0.00
64) Phenanthrene-d10	17.31	188	566613	20.00	ng	0.00
76) Chrysene-d12	21.55	240	533038	20.00	ng	0.00
86) Perylene-d12	24.66	264	558889	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	472761	92.33	ng	0.00
7) Phenol-d6	7.10	99	657939	90.65	ng	0.00
23) Nitrobenzene-d5	9.09	82	433152	64.18	ng	0.00
42) 2,4,6-Tribromophenol	16.05	330	291588	81.74	ng	0.00
45) 2-Fluorobiphenyl	13.18	172	1159566	67.23	ng	0.00
79) Terphenyl-d14	19.92	244	1481155	59.14	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.01	163	293097	14.389	ng	99
71) Phenanthrene	17.34	178	206579	7.193	ng	96
75) Fluoranthene	19.36	202	639246	17.296	ng	96
78) Pyrene	19.72	202	704338	20.750	ng	97
81) Benzo(a)anthracene	21.53	228	429046	12.914	ng	97
83) Chrysene	21.60	228	409118	12.801	ng	99
87) Indeno(1,2,3-cd)pyrene	28.17	276	307429	7.659	ng	97
88) Benzo(b)fluoranthene	23.68	252	582893	16.703	ng	98
89) Benzo(k)fluoranthene	23.73	252	175835m	5.213	ng	
90) Benzo(a)pyrene	24.51	252	445134	13.668	ng	99
91) Dibenzo(a,h)anthracene	28.21	278	77635m	2.397	ng	
92) Benzo(a,h,i)perylene	29.26	276	333192	10.301	ng	95

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

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