

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG083120\
 Data File : BG046477.D
 Acq On : 31 Aug 2020 20:32
 Operator : CG/JU
 Sample : L3847-17
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LINDEN-JOP-BH-10-C

Manual Integrations
 APPROVED

mohammad
 9/1/2020 10:47:32 AM

Quant Time: Sep 07 04:13:51 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	71397	20.00	ng	0.00
21) Naphthalene-d8	10.73	136	294956	20.00	ng	0.00
39) Acenaphthene-d10	14.56	164	213855	20.00	ng	0.00
64) Phenanthrene-d10	17.30	188	531373	20.00	ng	0.00
76) Chrysene-d12	21.55	240	528156	20.00	ng	0.00
86) Perylene-d12	24.65	264	561016	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	335593	83.25	ng	0.00
7) Phenol-d6	7.10	99	451458	79.00	ng	0.00
23) Nitrobenzene-d5	9.08	82	288561	55.92	ng	0.00
42) 2,4,6-Tribromophenol	16.05	330	222663	76.84	ng	0.00
45) 2-Fluorobiphenyl	13.18	172	800804	57.16	ng	0.00
79) Terphenyl-d14	19.92	244	1306462	52.64	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.01	163	33543	2.027	ng	95
71) Phenanthrene	17.35	178	590567	21.927	ng	98
72) Anthracene	17.44	178	119911	4.512	ng	96
75) Fluoranthene	19.36	202	763343	22.023	ng	98
78) Pyrene	19.72	202	685694	20.388	ng	98
81) Benzo(a)anthracene	21.53	228	340443	10.342	ng	97
83) Chrysene	21.60	228	315277m	9.956	ng	
87) Indeno(1,2,3-cd)pyrene	28.15	276	190961	4.739	ng	95
88) Benzo(b)fluoranthene	23.67	252	373348	10.658	ng	99
89) Benzo(k)fluoranthene	23.73	252	137372m	4.058	ng	
90) Benzo(a)pyrene	24.50	252	283801	8.681	ng	99
92) Benzo(a,h,i)perylene	29.24	276	188673	5.811	ng	96

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

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