

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF083120\
 Data File : BF121489.D
 Acq On : 31 Aug 2020 21:04
 Operator : JU/CG
 Sample : L3847-29
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LINDEN-JOP-BH-12-C

Manual Integrations
 APPROVED

mohammad
 9/2/2020 1:56:20 PM

Quant Time: Sep 10 12:33:29 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 28 09:25:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	294338	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	1029341	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	495263	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	796357	20.00	ng	0.00
76) Chrysene-d12	14.03	240	620817	20.00	ng	0.00
86) Perylene-d12	15.50	264	601813	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	1464415	83.47	ng	0.00
7) Phenol-d6	6.49	99	1892507	77.19	ng	0.00
23) Nitrobenzene-d5	7.42	82	1197919	79.56	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	434928	86.94	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	1959400	63.84	ng	0.00
79) Terphenyl-d14	12.97	244	1454992	49.33	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.62	163	527238	15.498	ng	98
71) Phenanthrene	11.41	178	331943	8.122	ng	96
75) Fluoranthene	12.60	202	435666	9.891	ng	97
78) Pyrene	12.83	202	522352	11.781	ng	96
81) Benzo(a)anthracene	14.02	228	287847	7.595	ng	96
83) Chrysene	14.05	228	281439	7.423	ng	95
87) Indeno(1,2,3-cd)pyrene	16.97	276	167504	4.524	ng	98
88) Benzo(b)fluoranthene	15.07	252	354777m	9.063	ng	
90) Benzo(a)pyrene	15.44	252	319759	9.084	ng	# 94
92) Benzo(g,h,i)perylene	17.42	276	245520	8.409	ng	97

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

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