

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083018\
 Data File : BF108650.D
 Acq On : 30 Aug 2018 19:50
 Operator : JU/SJ
 Sample : J4275-21 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-082918-B

Manual Integrations
 APPROVED

Sohil
 8/31/2018 12:18:19 PM

Quant Time: Aug 31 09:16:58 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 30 15:06:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	85889	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	367380	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	168887	20.00	ng	0.00
63) Phenanthrene-d10	11.54	188	288995	20.00	ng	-0.01
75) Chrysene-d12	14.19	240	283112	20.00	ng	0.00
86) Perylene-d12	15.75	264	282751	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.67	112	152399	30.82	ng	0.02
7) Phenol-d6	6.65	99	286232	40.42	ng	0.00
23) Nitrobenzene-d5	7.57	82	191779	26.50	ng	0.00
41) 2,4,6-Tribromophenol	10.84	330	72714	32.61	ng	-0.01
44) 2-Fluorobiphenyl	9.35	172	367500	29.26	ng	0.00
78) Terphenyl-d14	13.12	244	311962	21.41	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.75	163	33569	2.524	ng	# 97
74) Fluoranthene	12.76	202	41656	2.537	ng	95
80) Benzo(a)anthracene	14.18	228	44213	2.562	ng	# 85
82) Chrysene	14.22	228	54544m	3.286	ng	
85) Indeno(1,2,3-cd)pyrene	17.37	276	54088	4.687	ng	# 86
87) Benzo(b)fluoranthene	15.28	252	108619m	6.254	ng	
88) Benzo(k)fluoranthene	15.31	252	41669m	2.422	ng	
89) Benzo(a)pyrene	15.68	252	72100	4.593	ng	96
91) Benzo(g,h,i)perylene	17.86	276	47708	3.916	ng	# 88

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

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