

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082118\
 Data File : BF108353.D
 Acq On : 21 Aug 2018 22:30
 Operator : JU/SJ
 Sample : J4497-01RE
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CARLSTADT-FRONTRE

Manual Integrations
 APPROVED

Sohil
 8/22/2018 4:03:12 PM

Quant Time: Aug 22 00:44:16 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 21 15:20:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.01	152	176109	20.00	ng	0.00
21) Naphthalene-d8	8.29	136	641768	20.00	ng	-0.01
38) Acenaphthene-d10	10.05	164	292198	20.00	ng	-0.02
63) Phenanthrene-d10	11.54	188	491326	20.00	ng	-0.02
75) Chrysene-d12	14.19	240	375787	20.00	ng	-0.02
86) Perylene-d12	15.75	264	335014	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	940656	96.70	ng	0.00
7) Phenol-d6	6.63	99	1257101	97.25	ng	0.00
23) Nitrobenzene-d5	7.57	82	830204	72.19	ng	-0.01
41) 2,4,6-Tribromophenol	10.84	330	291576	100.04	ng	-0.02
44) 2-Fluorobiphenyl	9.36	172	1491202	81.93	ng	-0.02
78) Terphenyl-d14	13.13	244	1180083	79.84	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.75	163	108155	5.314	ng	100
70) Phenanthrene	11.57	178	75412	3.254	ng	98
74) Fluoranthene	12.76	202	151394	5.986	ng	96
77) Pyrene	13.00	202	141611	5.952	ng	98
80) Benzo(a)anthracene	14.18	228	76837	3.563	ng	98
82) Chrysene	14.22	228	72952	3.690	ng	97
85) Indeno(1,2,3-cd)pyrene	17.36	276	41166	2.403	ng	# 85
87) Benzo(b)fluoranthene	15.28	252	100927m	5.042	ng	
89) Benzo(a)pyrene	15.68	252	63424	3.538	ng	97
91) Benzo(g,h,i)perylene	17.85	276	39276	2.689	ng	# 90

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

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