

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
 Data File : BF110845.D
 Acq On : 19 Nov 2018 11:44
 Operator : JU/SJ
 Sample : J5951-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FDSB-02(21-23)

Manual Integrations
 APPROVED

Sohil
 11/20/2018 11:07:15 AM

Quant Time: Nov 19 13:14:27 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	69983	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	303448	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	146178	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	274685	20.00	ng	-0.01
76) Chrysene-d12	14.13	240	166009	20.00	ng	-0.01
87) Perylene-d12	15.66	264	157396	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	404412	93.86	ng	0.00
7) Phenol-d6	6.57	99	512403	93.00	ng	0.00
23) Nitrobenzene-d5	7.50	82	325722	61.82	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	133639	90.73	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	606252	59.23	ng	-0.01
79) Terphenyl-d14	13.06	244	538258	60.72	ng	-0.01
Target Compounds						Qvalue
50) Dimethylphthalate	9.69	163	54763	5.021	ng	99
71) Phenanthrene	11.50	178	136969	9.307	ng	99
72) Anthracene	11.56	178	44134	2.973	ng	98
75) Fluoranthene	12.69	202	106193	7.198	ng	98
78) Pyrene	12.93	202	83937	6.567	ng	97
81) Benzo(a)anthracene	14.12	228	28090	2.831	ng	97
83) Chrysene	14.16	228	24683	2.611	ng	98
88) Benzo(b)fluoranthene	15.20	252	26410m	2.805	ng	
90) Benzo(a)pyrene	15.59	252	20406	2.385	ng	99

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

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