

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070918\
 Data File : BF107235.D
 Acq On : 9 Jul 2018 18:08
 Operator : JU/SJ
 Sample : J3880-11RE
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SURCHARGE-SP-16RE

Manual Integrations
 APPROVED

Sohil
 7/10/2018 2:46:37 PM

Quant Time: Jul 10 04:16:12 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	32850	20.00	ng	-0.01
21) Naphthalene-d8	8.23	136	113440	20.00	ng	-0.02
38) Acenaphthene-d10	9.99	164	48908	20.00	ng	-0.02
63) Phenanthrene-d10	11.49	188	99491	20.00	ng	-0.01
75) Chrysene-d12	14.15	240	90519	20.00	ng	-0.02
86) Perylene-d12	15.69	264	57358	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	188826	97.33	ng	0.00
7) Phenol-d6	6.60	99	228095	94.15	ng	-0.02
23) Nitrobenzene-d5	7.51	82	129748	63.56	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	54784	96.65	ng	-0.01
44) 2-Fluorobiphenyl	9.30	172	230102	78.51	ng	-0.02
78) Terphenyl-d14	13.07	244	303789	81.29	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.70	163	56721	15.463	ng	98
70) Phenanthrene	11.51	178	31234	6.509	ng	99
71) Anthracene	11.57	178	13035	2.661	ng	99
74) Fluoranthene	12.71	202	87335	16.512	ng	93
77) Pyrene	12.94	202	87153	14.601	ng	98
80) Benzo(a)anthracene	14.14	228	43174m	7.826	ng	
82) Chrysene	14.17	228	44514m	8.770	ng	
85) Indeno(1,2,3-cd)pyrene	17.29	276	18959	4.402	ng	98
87) Benzo(b)fluoranthene	15.23	252	41657m	11.691	ng	
88) Benzo(k)fluoranthene	15.26	252	14702m	4.333	ng	
89) Benzo(a)pyrene	15.62	252	28050	8.856	ng	# 91
90) Dibenzo(a,h)anthracene	17.28	278	5614	2.091	ng	# 68
91) Benzo(a,h,i)perylene	17.78	276	22403	8.538	ng	# 87

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

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