

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100318\
 Data File : BF109555.D
 Acq On : 3 Oct 2018 19:10
 Operator : JU/SJ
 Sample : J5198-09 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-100218-E

Manual Integrations
 APPROVED

Sohil
 10/4/2018 11:44:26 AM

Quant Time: Oct 04 02:27:35 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	75525	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	264560	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	112695	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	236766	20.00	ng	0.00
75) Chrysene-d12	13.84	240	220391	20.00	ng	0.00
86) Perylene-d12	15.24	264	151906	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	174018	37.95	ng	0.00
7) Phenol-d6	6.35	99	234051	40.00	ng	-0.01
23) Nitrobenzene-d5	7.26	82	132826	29.64	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	53919	48.53	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	225578	30.19	ng	-0.02
78) Terphenyl-d14	12.79	244	261628	29.13	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.45	163	77760	9.487	ng	Qvalue # 98
51) Acenaphthene	9.76	154	18003	2.691	ng	96
57) Fluorene	10.28	166	29216	4.115	ng	99
70) Phenanthrene	11.23	178	136719	11.565	ng	98
71) Anthracene	11.29	178	51357	4.283	ng	100
74) Fluoranthene	12.42	202	211670	16.755	ng	95
77) Pyrene	12.64	202	196790	12.459	ng	100
80) Benzo(a)anthracene	13.83	228	83052	6.256	ng	97
82) Chrysene	13.87	228	77670	5.903	ng	97
85) Indeno(1,2,3-cd)pyrene	16.60	276	24744	2.320	ng	# 91
87) Benzo(b)fluoranthene	14.85	252	72202m	8.650	ng	
88) Benzo(k)fluoranthene	14.87	252	19593m	2.474	ng	
89) Benzo(a)pyrene	15.19	252	47883	6.366	ng	# 91
91) Benzo(a,h,i)perylene	17.01	276	24388	4.021	ng	# 89

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

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