

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061918\
 Data File : BF106618.D
 Acq On : 20 Jun 2018 9:02
 Operator : JU/SJ
 Sample : J3249-05DL 25X
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-01-061818-CDL

Manual Integrations
 APPROVED

Sohil
 6/20/2018 4:58:51 PM

Quant Time: Jun 20 09:25:15 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 19 14:32:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	91012	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	320831	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	127703	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	223807	20.00	ng	0.00
75) Chrysene-d12	14.15	240	254934	20.00	ng	0.00
86) Perylene-d12	15.69	264	185452	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	21649	4.03	ng	0.00
7) Phenol-d6	6.60	99	28143	4.19	ng	-0.01
23) Nitrobenzene-d5	7.52	82	15300	2.65	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	4637	3.13	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	31968	-12.72	ng	-0.01
78) Terphenyl-d14	13.08	244	34746	3.30	ng	0.00
Target Compounds						
51) Acenaphthene	10.04	154	17467	2.193	ng	Qvalue 100
57) Fluorene	10.55	166	34062	3.936	ng	98
70) Phenanthrene	11.52	178	299188	27.716	ng	99
71) Anthracene	11.57	178	103393	9.384	ng	99
74) Fluoranthene	12.72	202	545624	45.859	ng	96
77) Pyrene	12.95	202	490455	29.174	ng	100
80) Benzo(a)anthracene	14.14	228	295695m	19.031	ng	
82) Chrysene	14.18	228	257162m	17.990	ng	
85) Indeno(1,2,3-cd)pyrene	17.25	276	86971	7.170	ng	# 92
87) Benzo(b)fluoranthene	15.24	252	256671m	22.280	ng	
88) Benzo(k)fluoranthene	15.26	252	93815m	8.551	ng	
89) Benzo(a)pyrene	15.62	252	175071	17.095	ng	96
90) Dibenzo(a,h)anthracene	17.26	278	22417	2.583	ng	93
91) Benzo(g,h,i)perylene	17.74	276	86071	10.146	ng	# 92

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

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