

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062518\
 Data File : BF106763.D
 Acq On : 25 Jun 2018 11:32
 Operator : JU/SJ
 Sample : J3246-03DL 4X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-062018DL

Manual Integrations
 APPROVED

Sohil
 6/27/2018 10:55:30 AM

Quant Time: Jun 25 14:05:07 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 22 11:34:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	49727	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	202054	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	104323	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	200167	20.00	ng	0.00
75) Chrysene-d12	14.15	240	160418	20.00	ng	0.00
86) Perylene-d12	15.70	264	172130	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	72219	24.59	ng	0.00
7) Phenol-d6	6.60	99	93849	25.59	ng	-0.01
23) Nitrobenzene-d5	7.52	82	55789	15.34	ng	-0.01
41) 2,4,6-Tribromophenol	10.79	330	29863	24.70	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	129029	7.48	ng	-0.01
78) Terphenyl-d14	13.07	244	136864	20.67	ng	-0.01
Target Compounds						
31) Naphthalene	8.26	128	40095	4.244	ng	98
37) 2-Methylnaphthalene	8.95	142	29535	4.717	ng	92
51) Acenaphthene	10.03	154	44622	6.858	ng	97
54) Dibenzofuran	10.20	168	42954	4.821	ng	94
57) Fluorene	10.55	166	72898	10.311	ng	99
70) Phenanthrene	11.52	178	520545	53.918	ng	99
71) Anthracene	11.57	178	128847	13.075	ng	99
72) Carbazole	11.72	167	74317	8.055	ng	98
74) Fluoranthene	12.71	202	521414	49.000	ng	96
77) Pyrene	12.95	202	456243	43.129	ng	100
80) Benzo(a)anthracene	14.14	228	216221	22.115	ng	99
82) Chrysene	14.18	228	202796	22.546	ng	97
85) Indeno(1,2,3-cd)pyrene	17.27	276	105060	13.763	ng	# 87
87) Benzo(b)fluoranthene	15.24	252	259586m	24.277	ng	
88) Benzo(k)fluoranthene	15.27	252	91124m	8.949	ng	
89) Benzo(a)pyrene	15.64	252	194933m	20.507	ng	
90) Dibenzo(a,h)anthracene	17.28	278	25401	3.153	ng	# 92
91) Benzo(g,h,i)perylene	17.75	276	96772	12.290	ng	# 89

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

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