

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
 Data File : BF103211.D
 Acq On : 26 Feb 2018 11:54
 Operator : SJ/JU
 Sample : J1398-05 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-022218

Manual Integrations
 APPROVED

Sohil
 2/27/2018 12:22:28 PM

Quant Time: Feb 26 13:23:46 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 23 02:43:46 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	159336	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	653814	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	264783	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	381862	20.00	ng	0.00
75) Chrysene-d12	14.04	240	280837	20.00	ng	0.00
86) Perylene-d12	15.53	264	299393	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	572242	54.32	ng	0.00
7) Phenol-d6	6.50	99	670204	51.99	ng	0.00
23) Nitrobenzene-d5	7.43	82	430367	37.59	ng	-0.01
41) 2,4,6-Tribromophenol	10.70	330	95893	42.79	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	597484	32.89	ng	0.00
78) Terphenyl-d14	12.97	244	381701	32.67	ng	-0.01
Target Compounds						Qvalue
51) Acenaphthene	9.94	154	37729	2.396	ng	98
57) Fluorene	10.46	166	49439	2.685	ng	100
70) Phenanthrene	11.42	178	324786	15.455	ng	99
71) Anthracene	11.47	178	123663	5.739	ng	97
74) Fluoranthene	12.61	202	531583	25.172	ng	99
77) Pyrene	12.84	202	464326	21.206	ng	99
80) Benzo(a)anthracene	14.03	228	250588	13.752	ng	98
82) Chrysene	14.07	228	235443	13.307	ng	98
85) Indeno(1,2,3-cd)pyrene	17.04	276	121245	8.656	ng	96
87) Benzo(b)fluoranthene	15.10	252	322618m	16.389	ng	
88) Benzo(k)fluoranthene	15.12	252	108261m	5.840	ng	
89) Benzo(a)pyrene	15.47	252	232987	13.254	ng	# 93
90) Dibenzo(a,h)anthracene	17.04	278	29335	2.160	ng	# 80
91) Benzo(g,h,i)perylene	17.50	276	119386	8.993	ng	95

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

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