

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF032018\
 Data File : BF103814.D
 Acq On : 20 Mar 2018 20:36
 Operator : JU/SJ
 Sample : J1982-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-111(5-6)

Manual Integrations
 APPROVED

Sohil
 3/22/2018 12:57:46 PM

Quant Time: Mar 21 12:44:20 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 15 18:31:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	147652	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	557505	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	235454	20.00	ng	0.01
63) Phenanthrene-d10	11.51	188	432882	20.00	ng	0.01
75) Chrysene-d12	14.16	240	275885	20.00	ng	0.01
86) Perylene-d12	15.70	264	251432	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	1316781	135.64	ng	0.02
7) Phenol-d6	6.60	99	1632108	137.42	ng	0.02
23) Nitrobenzene-d5	7.54	82	874819	109.43	ng	0.00
41) 2,4,6-Tribromophenol	10.82	330	315466	155.83	ng	0.02
44) 2-Fluorobiphenyl	9.33	172	1217407	82.48	ng	0.00
78) Terphenyl-d14	13.10	244	1018056	85.65	ng	0.01
Target Compounds						
31) Naphthalene	8.28	128	309597	10.993	ng	# 96
48) Acenaphthylene	9.88	152	110242	4.559	ng	# 84
49) Dimethylphthalate	9.72	163	141801	7.894	ng	# 95
51) Acenaphthene	10.06	154	347390	24.194	ng	98
54) Dibenzofuran	10.22	168	311724	15.201	ng	# 90
57) Fluorene	10.57	166	656130	41.233	ng	99
70) Phenanthrene	11.53	178	635801	26.850	ng	100
71) Anthracene	11.59	178	800179	33.271	ng	100
72) Carbazole	11.74	167	261246	11.176	ng	99
74) Fluoranthene	12.74	202	841246	34.484	ng	100
77) Pyrene	12.96	202	693747	34.241	ng	99
80) Benzo(a)anthracene	14.15	228	317778	18.197	ng	97
82) Chrysene	14.19	228	241810m	14.526	ng	
85) Indeno(1,2,3-cd)pyrene	17.27	276	88111	6.061	ng	95
87) Benzo(b)fluoranthene	15.24	252	293893m	18.501	ng	
88) Benzo(k)fluoranthene	15.27	252	84693m	5.547	ng	
89) Benzo(a)pyrene	15.63	252	210780	14.630	ng	97
91) Benzo(g,h,i)perylene	17.75	276	82219	6.774	ng	96

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

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