

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF022818\
 Data File : BF103298.D
 Acq On : 28 Feb 2018 17:37
 Operator : SJ/JU
 Sample : J1399-11
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-03-022718-B

Manual Integrations
 APPROVED

Sohil
 3/1/2018 11:43:04 AM

Quant Time: Mar 01 00:34:00 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 17:05:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	141581	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	557787	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	213185	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	285302	20.00	ng	0.00
75) Chrysene-d12	14.06	240	230096	20.00	ng	0.00
86) Perylene-d12	15.56	264	220287	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	744446	79.52	ng	0.00
7) Phenol-d6	6.50	99	908416	79.30	ng	0.00
23) Nitrobenzene-d5	7.43	82	602687	61.71	ng	-0.01
41) 2,4,6-Tribromophenol	10.70	330	113797	63.06	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	812345	61.34	ng	0.00
78) Terphenyl-d14	12.99	244	506458	52.91	ng	0.00
Target Compounds						
31) Naphthalene	8.17	128	266940	9.775	ng	100
37) 2-Methylnaphthalene	8.87	142	111346	6.309	ng	98
49) Dimethylphthalate	9.62	163	74270	4.343	ng	100
51) Acenaphthene	9.95	154	79741	6.289	ng	100
54) Dibenzofuran	10.12	168	69836	4.012	ng	97
57) Fluorene	10.46	166	88558	5.973	ng	99
70) Phenanthrene	11.43	178	464622	29.592	ng	99
71) Anthracene	11.48	178	140543	8.729	ng	98
72) Carbazole	11.64	167	72753	4.538	ng	99
74) Fluoranthene	12.62	202	469457	29.754	ng	98
77) Pyrene	12.86	202	411449	22.935	ng	99
80) Benzo(a)anthracene	14.04	228	203200m	13.611	ng	
82) Chrysene	14.08	228	194293m	13.403	ng	
85) Indeno(1,2,3-cd)pyrene	17.08	276	71398	6.221	ng	96
87) Benzo(b)fluoranthene	15.12	252	215542m	14.882	ng	
88) Benzo(k)fluoranthene	15.14	252	79115m	5.801	ng	
89) Benzo(a)pyrene	15.49	252	154358	11.934	ng	# 92
91) Benzo(g,h,i)perylene	17.55	276	66387	6.797	ng	# 92

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

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