

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF022818\
 Data File : BF103306.D
 Acq On : 28 Feb 2018 21:11
 Operator : SJ/JU
 Sample : J1716-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-01-08-12

Manual Integrations
 APPROVED

Sohil
 3/1/2018 11:43:16 AM

Quant Time: Mar 01 01:04:01 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	157238	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	685723	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	278083	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	397791	20.00	ng	0.00
75) Chrysene-d12	14.05	240	258968	20.00	ng	0.00
86) Perylene-d12	15.55	264	269798	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.50	112	1349444	129.80	ng	0.01
7) Phenol-d6	6.52	99	1641324	129.02	ng	0.00
23) Nitrobenzene-d5	7.44	82	1081463	90.07	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	270177	114.78	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1328854	81.51	ng	0.00
78) Terphenyl-d14	12.99	244	850976	78.99	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
31) Naphthalene	8.17	128	318873	9.498	ng	99
49) Dimethylphthalate	9.62	163	200564	8.991	ng	99
70) Phenanthrene	11.43	178	130425	5.958	ng	98
74) Fluoranthene	12.62	202	217812	9.901	ng	99
77) Pyrene	12.85	202	222044	10.997	ng	99
80) Benzo(a)anthracene	14.04	228	136903	8.148	ng	99
82) Chrysene	14.08	228	128115	7.853	ng	99
85) Indeno(1,2,3-cd)pyrene	17.07	276	179882	13.927	ng	97
87) Benzo(b)fluoranthene	15.12	252	334258m	18.843	ng	
88) Benzo(k)fluoranthene	15.14	252	83020m	4.970	ng	
89) Benzo(a)pyrene	15.49	252	256496	16.192	ng	96
90) Dibenzo(a,h)anthracene	17.07	278	35222	2.877	ng	# 77
91) Benzo(a,h,i)perylene	17.54	276	186062	15.553	ng	97

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

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