

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022718\
 Data File : BF103278.D
 Acq On : 27 Feb 2018 22:01
 Operator : SJ/JU
 Sample : J1702-03
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-SF-03-003-B

Manual Integrations
 APPROVED

Sohil
 2/28/2018 11:18:04 AM

Quant Time: Feb 28 02:20:24 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	173816	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	688355	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	240499	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	333506	20.00	ng	0.00
75) Chrysene-d12	14.06	240	258559	20.00	ng	0.00
86) Perylene-d12	15.56	264	171507	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	844875	73.52	ng	0.00
7) Phenol-d6	6.51	99	1022233	72.69	ng	0.00
23) Nitrobenzene-d5	7.44	82	657956	54.59	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	112067	55.05	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	863806	57.07	ng	0.00
78) Terphenyl-d14	12.99	244	563167	52.36	ng	0.00
Target Compounds						
70) Phenanthrene	11.43	178	102182	5.567	ng	97
74) Fluoranthene	12.62	202	137401	7.450	ng	97
77) Pyrene	12.86	202	130106	6.454	ng	98
80) Benzo(a)anthracene	14.04	228	55373m	3.301	ng	
82) Chrysene	14.08	228	56102m	3.444	ng	
87) Benzo(b)fluoranthene	15.12	252	45306m	4.018	ng	
89) Benzo(a)pyrene	15.49	252	30058	2.985	ng	# 87
91) Benzo(g,h,i)perylene	17.54	276	19505	2.565	ng	# 90

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

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