

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF021618\
 Data File : BF103057.D
 Acq On : 16 Feb 2018 17:26
 Operator : SJ/JU
 Sample : J1399-07 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-02-021518-B

Manual Integrations
 APPROVED

Sohil
 2/19/2018 11:09:40 AM

Quant Time: Feb 16 23:27:28 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 16 14:34:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	171546	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	681151	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	267873	20.00	ng	-0.01
63) Phenanthrene-d10	11.40	188	399461	20.00	ng	0.00
75) Chrysene-d12	14.04	240	307553	20.00	ng	-0.01
86) Perylene-d12	15.54	264	242860	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	247400	21.90	ng	0.00
7) Phenol-d6	6.50	99	308603	22.69	ng	-0.01
23) Nitrobenzene-d5	7.43	82	173844	17.70	ng	-0.01
41) 2,4,6-Tribromophenol	10.70	330	42201	19.57	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	291691	18.07	ng	-0.01
78) Terphenyl-d14	12.98	244	210268	16.19	ng	-0.01
Target Compounds						
70) Phenanthrene	11.43	178	112357	5.050	ng	Qvalue 100
74) Fluoranthene	12.62	202	182747	8.164	ng	99
77) Pyrene	12.85	202	161059	6.678	ng	99
80) Benzo(a)anthracene	14.03	228	78510	4.059	ng	98
82) Chrysene	14.07	228	71368m	3.660	ng	
87) Benzo(b)fluoranthene	15.10	252	71730m	4.556	ng	
89) Benzo(a)pyrene	15.47	252	52671	3.716	ng	# 85
91) Benzo(g,h,i)perylene	17.50	276	27371	2.431	ng	# 90

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

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