

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030918\
 Data File : BF103556.D
 Acq On : 9 Mar 2018 14:52
 Operator : SJ/JU
 Sample : J1755-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-03-030818

Manual Integrations
 APPROVED

Sohil
 3/12/2018 3:50:11 PM

Quant Time: Mar 09 16:26:23 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 09 11:47:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	135595	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	564673	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	236043	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	351209	20.00	ng	0.00
75) Chrysene-d12	14.06	240	227564	20.00	ng	0.00
86) Perylene-d12	15.57	264	201981	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	994099	110.88	ng	0.00
7) Phenol-d6	6.50	99	1237485	112.80	ng	0.00
23) Nitrobenzene-d5	7.43	82	804691	81.38	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	160155	80.16	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	984649	68.59	ng	0.00
78) Terphenyl-d14	12.98	244	616754	65.15	ng	0.00
Target Compounds						
48) Acenaphthylene	9.77	152	60059	2.495	ng	98
70) Phenanthrene	11.43	178	82302	4.258	ng	99
73) Di-n-butylphthalate	11.94	149	335799	13.597	ng	99
74) Fluoranthene	12.62	202	100065	5.152	ng	98
77) Pyrene	12.85	202	146674	8.267	ng	98
80) Benzo(a)anthracene	14.04	228	55752m	3.776	ng	
82) Chrysene	14.09	228	61592	4.296	ng	95
87) Benzo(b)fluoranthene	15.13	252	42896m	3.230	ng	
89) Benzo(a)pyrene	15.51	252	42548	3.588	ng	# 90
91) Benzo(g,h,i)perylene	17.60	276	26176	2.923	ng	# 89

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

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