

Data Path : S:\HPCHEM1\BNA_F\DATA\BF031016\
 Data File : BF085343.D
 Acq On : 10 Mar 2016 19:59
 Operator : UM/SJ
 Sample : H1753-22
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-COMP4

Manual Integrations
 APPROVED

Sohil
 3/11/2016 4:58:36 PM

Quant Time: Mar 11 10:07:34 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 11 00:01:03 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	156851	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	551101	20.00	ng	-0.01
38) Acenaphthene-d10	9.98	164	242934	20.00	ng	0.00
63) Phenanthrene-d10	11.46	188	340582	20.00	ng	0.00
75) Chrysene-d12	14.11	240	248661	20.00	ng	0.00
86) Perylene-d12	15.57	264	253647	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	1194712	127.76	ng	0.01
7) Phenol-d6	6.57	99	1785821	145.77	ng	0.01
23) Nitrobenzene-d5	7.50	82	1007875	97.86	ng	-0.01
41) 2,4,6-Tribromophenol	10.77	330	259302	124.74	ng	-0.01
44) 2-Fluorobiphenyl	9.30	172	1567891	98.15	ng	0.00
78) Terphenyl-d14	13.05	244	989908	92.41	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.69	163	236654	12.69	ng	99
70) Phenanthrene	11.49	178	206773	11.58	ng	97
71) Anthracene	11.53	178	57573	3.04	ng	99
74) Fluoranthene	12.68	202	245166	13.69	ng	97
77) Pyrene	12.91	202	240877	12.87	ng	97
80) Benzo(a)anthracene	14.09	228	112348m	7.55	ng	
82) Chrysene	14.13	228	98881m	7.19	ng	
85) Indeno(1,2,3-cd)pyrene	17.02	276	52824	4.92	ng	# 90
87) Benzo(b)fluoranthene	15.15	252	135201m	8.00	ng	
88) Benzo(k)fluoranthene	15.16	252	36984m	2.71	ng	
89) Benzo(a)pyrene	15.50	252	94642	6.76	ng	# 90
91) Benzo(g,h,i)perylene	17.47	276	50683	4.21	ng	94

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

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