

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050316\
 Data File : BF086649.D
 Acq On : 3 May 2016 17:25
 Operator : UM/SJ
 Sample : H2685-07
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26WARD-2B-CS-4(0-10FTBGS)

Manual Integrations
 APPROVED

sohil
 5/4/2016 6:52:57 PM

Quant Time: May 04 04:08:42 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 27 14:20:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	196011	20.00	ng	-0.06
21) Naphthalene-d8	8.04	136	796413	20.00	ng	-0.06
38) Acenaphthene-d10	9.79	164	345921	20.00	ng	-0.06
63) Phenanthrene-d10	11.26	188	524960	20.00	ng	-0.06
75) Chrysene-d12	13.91	240	402654	20.00	ng	-0.06
86) Perylene-d12	15.30	264	355151	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	1043789	91.85	ng	-0.04
7) Phenol-d6	6.40	99	1420161	89.54	ng	-0.05
23) Nitrobenzene-d5	7.32	82	1011652	68.47	ng	-0.06
41) 2,4,6-Tribromophenol	10.58	330	290756	79.70	ng	-0.06
44) 2-Fluorobiphenyl	9.12	172	1337680	63.78	ng	-0.06
78) Terphenyl-d14	12.85	244	933794	51.09	ng	-0.06
Target Compounds						
49) Dimethylphthalate	9.52	163	167953	6.45	ng	Qvalue 98
70) Phenanthrene	11.29	178	144947	5.26	ng	99
74) Fluoranthene	12.48	202	288652	10.38	ng	97
77) Pyrene	12.70	202	263929	9.10	ng	99
80) Benzo(a)anthracene	13.89	228	115288	5.37	ng	99
82) Chrysene	13.93	228	133040m	5.70	ng	
83) Bis(2-ethylhexyl)phthalate	13.89	149	33474	2.15	ng	# 98
85) Indeno(1,2,3-cd)pyrene	16.63	276	74207	4.29	ng	93
87) Benzo(b)fluoranthene	14.91	252	178934m	8.56	ng	
88) Benzo(k)fluoranthene	14.93	252	62114m	2.85	ng	
89) Benzo(a)pyrene	15.24	252	114390	5.80	ng	97
91) Benzo(a,h,i)perylene	17.03	276	71377	4.41	ng	93

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

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