

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042816\
 Data File : BF086483.D
 Acq On : 29 Apr 2016 00:46
 Operator : UM/SJ
 Sample : H2654-12
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26WARD-2A-CS-6(0-6FTBGS)

Manual Integrations
 APPROVED

sohil
 4/29/2016 4:52:53 PM

Quant Time: Apr 29 02:08:21 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	6.78	152	145592	20.00	ng	-0.02
21) Naphthalene-d8	8.08	136	608044	20.00	ng	-0.02
38) Acenaphthene-d10	9.82	164	249995	20.00	ng	-0.02
63) Phenanthrene-d10	11.31	188	380543	20.00	ng	-0.01
75) Chrysene-d12	13.94	240	308662	20.00	ng	-0.02
86) Perylene-d12	15.35	264	278362	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	864319	102.40	ng	0.00
7) Phenol-d6	6.43	99	1201463	101.99	ng	-0.01
23) Nitrobenzene-d5	7.36	82	800062	70.92	ng	-0.02
41) 2,4,6-Tribromophenol	10.61	330	193492	73.39	ng	-0.02
44) 2-Fluorobiphenyl	9.15	172	1075971	72.94	ng	-0.02
78) Terphenyl-d14	12.89	244	698050	49.82	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	9.55	163	231136	12.28	ng	99
70) Phenanthrene	11.33	178	81542	4.08	ng	97
74) Fluoranthene	12.51	202	138634	6.88	ng	98
77) Pyrene	12.74	202	128821	5.79	ng	98
80) Benzo(a)anthracene	13.93	228	74142	4.51	ng	99
82) Chrysene	13.96	228	56463	3.16	ng	97
85) Indeno(1,2,3-cd)pyrene	16.68	276	29137	2.20	ng	97
87) Benzo(b)fluoranthene	14.95	252	65563m	4.00	ng	
88) Benzo(k)fluoranthene	14.96	252	34885m	2.04	ng	
89) Benzo(a)pyrene	15.28	252	50092	3.24	ng	# 88

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

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