

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031416\
 Data File : BF085449.D
 Acq On : 15 Mar 2016 2:04
 Operator : UM/SJ
 Sample : H1730-10
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W170TH-SB-07-COMP(0.5-4.0)

Manual Integrations
 APPROVED

Sohil
 3/15/2016 2:09:36 PM

Quant Time: Mar 15 04:15:35 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 04 00:54:58 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	173378	20.00	ng	-0.06
21) Naphthalene-d8	8.21	136	667427	20.00	ng	-0.06
38) Acenaphthene-d10	9.97	164	319137	20.00	ng	-0.06
63) Phenanthrene-d10	11.45	188	539315	20.00	ng	-0.06
75) Chrysene-d12	14.09	240	331415	20.00	ng	-0.06
86) Perylene-d12	15.55	264	287326	20.00	ng	-0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	1642833	158.94	ng	-0.03
7) Phenol-d6	6.56	99	1738657	128.39	ng	-0.05
23) Nitrobenzene-d5	7.49	82	1172894	94.04	ng	-0.06
41) 2,4,6-Tribromophenol	10.76	330	392737	143.82	ng	-0.06
44) 2-Fluorobiphenyl	9.29	172	1821863	86.82	ng	-0.06
78) Terphenyl-d14	13.04	244	1286384	90.11	ng	-0.05
Target Compounds						Qvalue
49) Dimethylphthalate	9.68	163	415884	16.98	ng	99
70) Phenanthrene	11.48	178	103090	3.65	ng	96
74) Fluoranthene	12.66	202	234372	8.26	ng	96
77) Pyrene	12.89	202	221810	8.89	ng	98
80) Benzo(a)anthracene	14.08	228	108221	5.45	ng	97
82) Chrysene	14.12	228	93470m	5.10	ng	
85) Indeno(1,2,3-cd)pyrene	16.99	276	41835	2.92	ng	# 92
87) Benzo(b)fluoranthene	15.12	252	103861m	5.42	ng	
88) Benzo(k)fluoranthene	15.13	252	43806m	2.84	ng	
89) Benzo(a)pyrene	15.49	252	76826	4.84	ng	# 96
91) Benzo(g,h,i)perylene	17.43	276	37332	2.74	ng	96

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

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