

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031416\
 Data File : BF085447.D
 Acq On : 15 Mar 2016 1:08
 Operator : UM/SJ
 Sample : H1730-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W170TH-SB-23-COMP(0.5-1.5)

Manual Integrations
 APPROVED

Sohil
 3/15/2016 2:09:32 PM

Quant Time: Mar 15 04:09:14 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	166584	20.00	ng	-0.06
21) Naphthalene-d8	8.21	136	649960	20.00	ng	-0.06
38) Acenaphthene-d10	9.97	164	285345	20.00	ng	-0.06
63) Phenanthrene-d10	11.45	188	449893	20.00	ng	-0.06
75) Chrysene-d12	14.08	240	275827	20.00	ng	-0.07
86) Perylene-d12	15.55	264	283271	20.00	ng	-0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	1132287	114.01	ng	-0.06
7) Phenol-d6	6.55	99	1495072	114.91	ng	-0.06
23) Nitrobenzene-d5	7.49	82	855323	70.42	ng	-0.06
41) 2,4,6-Tribromophenol	10.76	330	301359	123.43	ng	-0.06
44) 2-Fluorobiphenyl	9.28	172	1233854	65.76	ng	-0.07
78) Terphenyl-d14	13.03	244	764883	64.37	ng	-0.06
Target Compounds						Qvalue
49) Dimethylphthalate	9.68	163	330730	15.10	ng	99
70) Phenanthrene	11.48	178	80547	3.42	ng	96
74) Fluoranthene	12.65	202	165820	7.01	ng	95
77) Pyrene	12.88	202	147243	7.09	ng	98
80) Benzo(a)anthracene	14.07	228	85654	5.19	ng	97
82) Chrysene	14.11	228	61339m	4.02	ng	
85) Indeno(1,2,3-cd)pyrene	16.99	276	42267	3.55	ng	# 90
87) Benzo(b)fluoranthene	15.12	252	103589m	5.49	ng	
88) Benzo(k)fluoranthene	15.15	252	34804m	2.29	ng	
89) Benzo(a)pyrene	15.48	252	73424	4.69	ng	# 93
91) Benzo(g,h,i)perylene	17.42	276	40984	3.05	ng	97

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

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