

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031416\
 Data File : BF085423.D
 Acq On : 14 Mar 2016 12:32
 Operator : UM/SJ
 Sample : H1743-02DL 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MNR-HARMON-BLDG-30-RW1DL

Manual Integrations
 APPROVED

Sohil
 3/15/2016 2:09:04 PM

Quant Time: Mar 15 01:49:52 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	148891	20.00	ng	-0.06
21) Naphthalene-d8	8.21	136	599403	20.00	ng	-0.06
38) Acenaphthene-d10	9.97	164	306600	20.00	ng	-0.06
63) Phenanthrene-d10	11.45	188	531908	20.00	ng	-0.06
75) Chrysene-d12	14.09	240	330959	20.00	ng	-0.06
86) Perylene-d12	15.55	264	253887	20.00	ng	-0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	318848	35.92	ng	-0.06
7) Phenol-d6	6.54	99	248052	21.33	ng	-0.07
23) Nitrobenzene-d5	7.49	82	582830	52.03	ng	-0.06
41) 2,4,6-Tribromophenol	10.76	330	195512	74.52	ng	-0.06
44) 2-Fluorobiphenyl	9.29	172	953138	47.28	ng	-0.06
78) Terphenyl-d14	13.04	244	849908	59.61	ng	-0.05
Target Compounds						
20) 3+4-Methylphenols	7.33	107	40950	3.96	ng	# 80
31) Naphthalene	8.23	128	210526	7.14	ng	# 90
37) 2-Methylnaphthalene	8.93	142	1043290	55.16	ng	98
49) Dimethylphthalate	9.68	163	109570	4.66	ng	# 93
51) Acenaphthene	10.00	154	51824	2.75	ng	# 80
54) Dibenzofuran	10.17	168	62711	2.54	ng	# 78
57) Fluorene	10.52	166	98335	5.06	ng	# 97
70) Phenanthrene	11.48	178	133878	4.80	ng	99

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

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