

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060716\
 Data File : BF087800.D
 Acq On : 7 Jun 2016 22:20
 Operator : UM/SJ
 Sample : H3379-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-2

Manual Integrations
 APPROVED

sohil
 6/8/2016 7:13:51 PM

Quant Time: Jun 08 05:06:43 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 07 18:51:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	103189	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	440902	20.00	ng	0.00
38) Acenaphthene-d10	9.87	164	202732	20.00	ng	-0.01
63) Phenanthrene-d10	11.35	188	329329	20.00	ng	-0.01
75) Chrysene-d12	13.98	240	226560	20.00	ng	-0.01
86) Perylene-d12	15.39	264	220537	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	708567	117.39	ng	0.01
7) Phenol-d6	6.48	99	1023654	139.93	ng	0.00
23) Nitrobenzene-d5	7.40	82	576791	76.39	ng	-0.01
41) 2,4,6-Tribromophenol	10.66	330	229425	107.18	ng	-0.01
44) 2-Fluorobiphenyl	9.20	172	911595	69.64	ng	-0.01
78) Terphenyl-d14	12.94	244	688918	69.19	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.59	163	110842	7.55	ng	99
70) Phenanthrene	11.37	178	222678	12.89	ng	98
71) Anthracene	11.43	178	65008	3.73	ng	97
74) Fluoranthene	12.56	202	315309	17.29	ng	98
77) Pyrene	12.79	202	293750	17.47	ng	99
80) Benzo(a)anthracene	13.97	228	141287	10.94	ng	99
82) Chrysene	14.00	228	96615	7.95	ng	96
85) Indeno(1,2,3-cd)pyrene	16.77	276	68056	6.03	ng	# 100
87) Benzo(b)fluoranthene	14.99	252	151438m	9.85	ng	
88) Benzo(k)fluoranthene	15.01	252	38527m	3.32	ng	
89) Benzo(a)pyrene	15.34	252	109072	8.77	ng	# 94
91) Benzo(a,h,i)perylene	17.18	276	67336	5.67	ng	97

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

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