

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060716\
 Data File : BF087803.D
 Acq On : 7 Jun 2016 23:48
 Operator : UM/SJ
 Sample : H3379-09
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
ClientSampled :
 SS-6

Manual Integrations
APPROVED
 sohil
 6/8/2016 7:13:16 PM

Quant Time: Jun 08 05:21:06 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	100400	20.00	ng	0.00	
21) Naphthalene-d8	8.12	136	408689	20.00	ng	0.00	
38) Acenaphthene-d10	9.87	164	171880	20.00	ng	-0.01	
63) Phenanthrene-d10	11.35	188	274332	20.00	ng	-0.01	
75) Chrysene-d12	13.98	240	205664	20.00	ng	-0.01	
86) Perylene-d12	15.39	264	206027	20.00	ng	0.00	
System Monitoring Compounds							
5) 2-Fluorophenol	5.44	112	547918	93.30	ng	0.01	
7) Phenol-d6	6.48	99	779777	109.56	ng	0.00	
23) Nitrobenzene-d5	7.40	82	455326	65.06	ng	-0.01	
41) 2,4,6-Tribromophenol	10.66	330	146854	80.92	ng	-0.01	
44) 2-Fluorobiphenyl	9.20	172	722625	64.72	ng	-0.01	
78) Terphenyl-d14	12.94	244	536672	59.38	ng	-0.01	
Target Compounds							Qvalue
49) Dimethylphthalate	9.60	163	103278	8.29	ng		98
70) Phenanthrene	11.37	178	52478	3.65	ng		99
74) Fluoranthene	12.56	202	72935	4.80	ng		95
77) Pyrene	12.79	202	62226	4.08	ng		98
80) Benzo(a)anthracene	13.97	228	34306	2.93	ng		97
82) Chrysene	14.00	228	23799m	2.16	ng		
87) Benzo(b)fluoranthene	14.99	252	39178m	2.73	ng		
89) Benzo(a)pyrene	15.34	252	27894	2.40	ng		99

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

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