

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060816\
 Data File : BF087812.D
 Acq On : 8 Jun 2016 16:22
 Operator : UM/SJ
 Sample : H3378-10
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-20

Manual Integrations
 APPROVED

sohil
 6/9/2016 7:59:25 PM

Quant Time: Jun 09 07:26:54 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	6.82	152	99097	20.00	ng	-0.02
21) Naphthalene-d8	8.11	136	435128	20.00	ng	-0.01
38) Acenaphthene-d10	9.87	164	185139	20.00	ng	-0.01
63) Phenanthrene-d10	11.35	188	292089	20.00	ng	-0.01
75) Chrysene-d12	13.99	240	208282	20.00	ng	0.00
86) Perylene-d12	15.42	264	205675	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	353235	60.94	ng	-0.01
7) Phenol-d6	6.46	99	477124	67.92	ng	-0.02
23) Nitrobenzene-d5	7.38	82	229526	30.80	ng	-0.03
41) 2,4,6-Tribromophenol	10.66	330	92226	47.18	ng	-0.01
44) 2-Fluorobiphenyl	9.19	172	404241	30.77	ng	-0.02
78) Terphenyl-d14	12.95	244	306730	33.51	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.59	163	127626	9.52	ng	99
70) Phenanthrene	11.37	178	40118	2.62	ng	99
74) Fluoranthene	12.56	202	68608	4.24	ng	98
77) Pyrene	12.79	202	58488	3.78	ng	99
80) Benzo(a)anthracene	13.98	228	31856	2.68	ng	97
82) Chrysene	14.01	228	23846m	2.14	ng	
87) Benzo(b)fluoranthene	15.01	252	35812m	2.50	ng	
89) Benzo(a)pyrene	15.36	252	26338	2.27	ng	97

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

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