

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110816\
 Data File : BF091100.D
 Acq On : 8 Nov 2016 23:34
 Operator : UM/SJ
 Sample : H5537-10
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 V001-BF014-02

Manual Integrations
 APPROVED

umangi
 11/9/2016 10:56:55 AM

Quant Time: Nov 09 02:46:55 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 08 12:59:29 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	223054	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	919036	20.00	ng	0.00
38) Acenaphthene-d10	9.70	164	452193	20.00	ng	0.00
63) Phenanthrene-d10	11.17	188	613399	20.00	ng	0.00
75) Chrysene-d12	13.80	240	425886	20.00	ng	-0.01
86) Perylene-d12	15.19	264	420016	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.25	112	1325771	98.16	ng	0.01
7) Phenol-d6	6.32	99	1746276	95.26	ng	0.00
23) Nitrobenzene-d5	7.23	82	1078796	64.03	ng	-0.01
41) 2,4,6-Tribromophenol	10.49	330	365505	89.41	ng	0.00
44) 2-Fluorobiphenyl	9.02	172	1590660	60.56	ng	-0.01
78) Terphenyl-d14	12.76	244	1268350	74.32	ng	0.00
Target Compounds						
10) Phenol	6.33	94	123499	6.09	ng	Qvalue # 76
49) Dimethylphthalate	9.42	163	106783	3.60	ng	99
74) Fluoranthene	12.39	202	107630	3.18	ng	91
77) Pyrene	12.60	202	97018	3.48	ng	96
80) Benzo(a)anthracene	13.79	228	57056	2.36	ng	98
87) Benzo(b)fluoranthene	14.81	252	72166m	2.53	ng	

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

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