

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110816\
 Data File : BF091097.D
 Acq On : 8 Nov 2016 22:12
 Operator : UM/SJ
 Sample : H5537-07
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 V001-BF014-01

Manual Integrations
 APPROVED

umangi
 11/9/2016 10:56:52 AM

Quant Time: Nov 09 02:36:49 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	186846	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	764466	20.00	ng	0.00
38) Acenaphthene-d10	9.70	164	391600	20.00	ng	0.00
63) Phenanthrene-d10	11.17	188	617441	20.00	ng	0.00
75) Chrysene-d12	13.80	240	390207	20.00	ng	-0.01
86) Perylene-d12	15.17	264	329566	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.25	112	1212127	107.14	ng	0.01
7) Phenol-d6	6.32	99	1569433	102.20	ng	0.00
23) Nitrobenzene-d5	7.23	82	1057081	75.43	ng	-0.01
41) 2,4,6-Tribromophenol	10.49	330	387257	109.39	ng	0.00
44) 2-Fluorobiphenyl	9.02	172	1509021	66.34	ng	-0.01
78) Terphenyl-d14	12.76	244	1437968	91.96	ng	0.00
Target Compounds						
10) Phenol	6.33	94	140518	8.27	ng	Qvalue # 64
49) Dimethylphthalate	9.42	163	90203	3.51	ng	99
70) Phenanthrene	11.20	178	75403	2.30	ng	99
74) Fluoranthene	12.39	202	130436	3.83	ng	90
77) Pyrene	12.60	202	119956	4.69	ng	97
80) Benzo(a)anthracene	13.79	228	54437	2.46	ng	100
87) Benzo(b)fluoranthene	14.81	252	56366m	2.52	ng	

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

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