

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110816\
 Data File : BF091101.D
 Acq On : 9 Nov 2016 00:02
 Operator : UM/SJ
 Sample : H5537-14
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 V001-BF015-01

Manual Integrations
 APPROVED

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

Quant Time: Nov 09 02:51:34 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	182892	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	725139	20.00	ng	0.00
38) Acenaphthene-d10	9.70	164	347714	20.00	ng	0.00
63) Phenanthrene-d10	11.17	188	486859	20.00	ng	0.00
75) Chrysene-d12	13.80	240	336794	20.00	ng	-0.01
86) Perylene-d12	15.19	264	318966	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.25	112	1232849	111.33	ng	0.01
7) Phenol-d6	6.32	99	1629010	108.38	ng	0.00
23) Nitrobenzene-d5	7.23	82	1005251	75.62	ng	-0.01
41) 2,4,6-Tribromophenol	10.49	330	348594	110.89	ng	0.00
44) 2-Fluorobiphenyl	9.02	172	1470439	72.81	ng	-0.01
78) Terphenyl-d14	12.76	244	1162035	86.10	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.33	94	122526	7.37	ng	# 74
49) Dimethylphthalate	9.42	163	349465	15.33	ng	98
70) Phenanthrene	11.20	178	77423	2.99	ng	97
74) Fluoranthene	12.39	202	124475	4.64	ng	90
77) Pyrene	12.60	202	109859	4.98	ng	96
80) Benzo(a)anthracene	13.79	228	60917	3.19	ng	98
82) Chrysene	13.83	228	45683m	2.42	ng	
87) Benzo(b)fluoranthene	14.81	252	77805m	3.60	ng	
89) Benzo(a)pyrene	15.13	252	48564	2.58	ng	# 92

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

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