

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102115\
 Data File : BF082440.D
 Acq On : 22 Oct 2015 9:33
 Operator : UM/IZ
 Sample : G4117-11
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AB-AC-08(0-2)

Quant Time: Oct 22 13:58:22 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 21 13:20:24 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	107967	20.00	ng	-0.01
21) Naphthalene-d8	8.06	136	436472	20.00	ng	-0.01
38) Acenaphthene-d10	9.82	164	204831	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	338849	20.00	ng	0.00
75) Chrysene-d12	13.98	240	269280	20.00	ng	-0.03
86) Perylene-d12	15.49	264	245766	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	703001	131.41	ng	0.01
7) Phenol-d6	6.42	99	948368	136.23	ng	0.00
23) Nitrobenzene-d5	7.35	82	562792	86.59	ng	0.00
41) 2,4,6-Tribromophenol	10.61	330	146638	76.34	ng	-0.01
44) 2-Fluorobiphenyl	9.14	172	1064199	86.16	ng	0.00
78) Terphenyl-d14	12.89	244	656851	64.64	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.54	163	211867	14.69	ng	# 97
70) Phenanthrene	11.33	178	71790	4.32	ng	96
74) Fluoranthene	12.51	202	85032	4.77	ng	99
77) Pyrene	12.74	202	78033	4.88	ng	98
80) Benzo(a)anthracene	13.97	228	37945	2.71	ng	97
82) Chrysene	14.01	228	35288	2.39	ng	97
87) Benzo(b)fluoranthene	15.08	252	45070m	3.01	ng	
89) Benzo(a)pyrene	15.43	252	26389	2.05	ng	# 95

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

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