

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090415\
 Data File : BE090724.D
 Acq On : 4 Sep 2015 11:33
 Operator : UM/IZ
 Sample : G3521-10
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X1SB0840006-20150901

Quant Time: Sep 04 13:20:07 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 02 11:40:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	42252	5.00	ng	0.00
6) Naphthalene-d8	10.31	136	187806	5.00	ng	0.00
12) Acenaphthene-d10	14.17	164	103464	5.00	ng	0.00
18) Phenanthrene-d10	16.91	188	259304	5.00	ng	0.00
25) Chrysene-d12	21.07	240	344098	5.00	ng	0.00
32) Perylene-d12	23.36	264	324641	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.18	112	88567	10.75	ng	0.00
5) Phenol-d6	6.74	99	127626	10.85	ng	0.00
7) Nitrobenzene-d5	8.71	82	56364	4.54	ng	0.00
13) 2,4,6-Tribromophenol	15.67	330	32192	10.55	ng	0.00
14) 2-Fluorobiphenyl	12.80	172	113524	3.96	ng	0.00
27) Terphenyl-d14	19.53	244	213659	5.61	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
22) Phenanthrene	16.96	178	2350	0.04	ng	# 80
24) Fluoranthene	18.96	202	5432	0.08	ng	# 96
26) Pyrene	19.32	202	5490	0.08	ng	# 96
28) Benzo(a)anthracene	21.05	228	4849	0.05	ng	# 95
31) Indeno(1,2,3-cd)pyrene	25.72	276	2644	0.03	ng	# 96
33) Benzo(b)fluoranthene	22.67	252	5744	0.08	ng	# 71
34) Benzo(k)fluoranthene	22.71	252	2296m	0.03	ng	# 71
35) Benzo(a)pyrene	23.26	252	3267	0.05	ng	# 65
37) Benzo(g,h,i)perylene	26.44	276	2858	0.04	ng	# 73

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

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