

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020516\
 Data File : BE091401.D
 Acq On : 5 Feb 2016 15:50
 Operator : UM/IZ
 Sample : H1339-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 OUTFALL001-160202

Quant Time: Feb 05 16:47:59 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE020216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 03 00:27:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	40443	5.00	ng	0.00
4) Naphthalene-d8	9.56	136	189950	5.00	ng	0.00
8) Acenaphthene-d10	13.35	164	124985	5.00	ng	0.00
14) Phenanthrene-d10	16.09	188	323113	5.00	ng	0.00
18) Chrysene-d12	20.22	240	385026	5.00	ng	0.00
24) Perylene-d12	22.10	264	370696	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	4.92	112	18569	2.10	ng	-0.02
3) Phenol-d6	6.58	99	13692	1.19	ng	0.02
5) Nitrobenzene-d5	8.17	82	41309	3.38	ng	0.00
9) 2,4,6-Tribromophenol	14.89	330	33810	5.63	ng	-0.01
10) 2-Fluorobiphenyl	12.01	172	145036	4.46	ng	0.00
20) Terphenyl-d14	18.61	244	290407	5.10	ng	0.00
Target Compounds						
7) 2-Methylnaphthalene	11.18	142	683	0.15	ng	Qvalue # 92
17) Fluoranthene	18.10	202	1781	0.14	ng	95
19) Pyrene	18.48	202	1823	0.02	ng	95
21) Benzo(a)anthracene	20.19	228	3156	0.21	ng	96
23) Indeno(1,2,3-cd)pyrene	23.73	276	4077	0.21	ng	96
25) Benzo(b)fluoranthene	21.56	252	2606	0.18	ng	# 81
26) Benzo(k)fluoranthene	21.58	252	3335	0.05	ng	# 84
27) Benzo(a)pyrene	22.01	252	2311	0.20	ng	# 70
28) Dibenzo(a,h)anthracene	23.68	278	4674	0.30	ng	# 92
29) Benzo(g,h,i)perylene	24.27	276	3799	0.18	ng	# 83

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

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