

Data Path : Z:\HPCHEM1\BNA E\DATA\BE030617\
 Data File : BE092797.D
 Acq On : 6 Mar 2017 21:25
 Operator : SJ/MA
 Sample : I2052-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 CPS093051

Manual Integrations
 APPROVED

mohammad
 3/7/2017 4:23:22 PM

Quant Time: Mar 07 07:35:40 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE030617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 06 15:46:01 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	9077	5.00	ng	-0.02
7) Naphthalene-d8	10.87	136	41216	5.00	ng	-0.02
12) Acenaphthene-d10	14.75	164	26132	5.00	ng	-0.02
18) Phenanthrene-d10	17.51	188	56992	5.00	ng	0.00
24) Chrysene-d12	21.68	240	83937	5.00	ng	-0.02
31) Perylene-d12	24.01	264	70893	5.00	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.60	112	8143	4.24	ng	-0.03
5) Phenol-d6	7.27	99	5994	2.62	ng	-0.05
8) Nitrobenzene-d5	9.27	82	12315	4.89	ng	-0.05
13) 2,4,6-Tribromophenol	16.24	330	9512	13.84	ng	-0.04
14) 2-Fluorobiphenyl	13.36	172	31635	5.13	ng	-0.04
26) Terphenyl-d14	20.10	244	50413	4.98	ng	-0.03
Target Compounds						
9) Naphthalene	10.91	128	8803	1.00	ng	Qvalue 94
11) 2-Methylnaphthalene	12.55	142	24370m	4.27	ng	
17) Fluorene	15.80	166	2115	0.28	ng	# 94
20) Pentachlorophenol	17.17	266	267005	313.07	ng	# 86
21) Phenanthrene	17.53	178	1259	0.09	ng	# 92
30) Indeno(1,2,3-cd)pyrene	26.48	276	2999	0.04	ng	97
36) Benzo(g,h,i)perylene	27.23	276	2881	0.04	ng	# 60

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

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