

Data Path : Z:\HPCHEM1\BNA E\DATA\BE011516\
 Data File : BE091372.D
 Acq On : 15 Jan 2016 9:38
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Quant Time: Jan 16 05:03:17 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE011416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 15 15:14:26 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	24505	5.00	ng	0.00
4) Naphthalene-d8	9.55	136	115297	5.00	ng	0.00
8) Acenaphthene-d10	13.36	164	58647	5.00	ng	0.00
14) Phenanthrene-d10	16.08	188	153229	5.00	ng	0.00
18) Chrysene-d12	20.21	240	171303	5.00	ng	0.00
24) Perylene-d12	22.10	264	145004	5.00	ng	0.00

System Monitoring Compounds						
2) 2-Fluorophenol	4.93	112	4130	1.50	ng	-0.01
3) Phenol-d6	6.56	99	5168	1.17	ng	0.00
5) Nitrobenzene-d5	8.16	82	6654	0.98	ng	0.00
9) 2,4,6-Tribromophenol	14.89	330	1169	1.03	ng	0.00
10) 2-Fluorobiphenyl	12.01	172	13692	1.03	ng	0.00
20) Terphenyl-d14	18.62	244	18626	1.03	ng	0.00

Target Compounds					Qvalue	
6) Naphthalene	9.59	128	24885	1.01	ng	100
7) 2-Methylnaphthalene	11.12	142	10418	1.01	ng	95
11) Acenaphthylene	13.10	152	83571	0.99	ng	100
12) Acenaphthene	13.42	154	16292	0.98	ng	83
13) Fluorene	14.40	166	19921	0.97	ng	97
15) Phenanthrene	16.11	178	39336	1.04	ng	100
16) Anthracene	16.20	178	30447	1.03	ng	100
17) Fluoranthene	18.10	202	31910	1.03	ng	99
19) Pyrene	18.48	202	27676	1.00	ng	99
21) Benzo(a)anthracene	20.19	228	26321	1.04	ng	99
22) Chrysene	20.25	228	42168	1.01	ng	99
23) Indeno(1,2,3-cd)pyrene	23.70	276	28185	0.98	ng	96
25) Benzo(b)fluoranthene	21.55	252	28966	1.03	ng	100
26) Benzo(k)fluoranthene	21.57	252	34388	1.05	ng	99
27) Benzo(a)pyrene	22.00	252	26520	0.97	ng	98
28) Dibenzo(a,h)anthracene	23.65	278	23065	1.08	ng	100
29) Benzo(g,h,i)perylene	24.25	276	27457	1.06	ng	98

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

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