

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020516\
 Data File : BE091402.D
 Acq On : 5 Feb 2016 16:29
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001EC

Quant Time: Feb 05 17:43:27 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	40576	5.00	ng	0.00
4) Naphthalene-d8	9.55	136	197397	5.00	ng	0.00
8) Acenaphthene-d10	13.35	164	128529	5.00	ng	0.00
14) Phenanthrene-d10	16.09	188	323452	5.00	ng	0.00
18) Chrysene-d12	20.22	240	391198	5.00	ng	0.00
24) Perylene-d12	22.09	264	374504	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	4.95	112	7295	0.96	ng	0.00
3) Phenol-d6	6.56	99	11548	1.03	ng	0.00
5) Nitrobenzene-d5	8.17	82	9998	1.00	ng	0.00
9) 2,4,6-Tribromophenol	14.90	330	3780	0.91	ng	0.00
10) 2-Fluorobiphenyl	12.01	172	31419	0.94	ng	0.00
20) Terphenyl-d14	18.62	244	56551	0.98	ng	0.00
Target Compounds						Qvalue
6) Naphthalene	9.59	128	41117	0.92	ng	99
7) 2-Methylnaphthalene	11.13	142	23554	0.96	ng	97
11) Acenaphthylene	13.10	152	178498	0.95	ng	100
12) Acenaphthene	13.42	154	31705	0.90	ng	98
13) Fluorene	14.40	166	42657	0.94	ng	99
15) Phenanthrene	16.12	178	78133	0.92	ng	100
16) Anthracene	16.21	178	69062	0.97	ng	99
17) Fluoranthene	18.10	202	77187	0.96	ng	99
19) Pyrene	18.48	202	81119	0.96	ng	100
21) Benzo(a)anthracene	20.19	228	69294	0.94	ng	98
22) Chrysene	20.25	228	95793	0.93	ng	98
23) Indeno(1,2,3-cd)pyrene	23.70	276	71934	0.94	ng	99
25) Benzo(b)fluoranthene	21.55	252	76889	0.95	ng	99
26) Benzo(k)fluoranthene	21.58	252	83987	0.91	ng	98
27) Benzo(a)pyrene	22.00	252	67943	0.93	ng	100
28) Dibenzo(a,h)anthracene	23.65	278	63485	0.96	ng	99
29) Benzo(a,h,i)perylene	24.25	276	76378	0.94	ng	98

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

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