

Data Path : Z:\HPCHEM1\BNA E\DATA\BE011516\
 Data File : BE091374.D
 Acq On : 15 Jan 2016 11:11
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001EC

Quant Time: Jan 16 05:04:15 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	22830	5.00	ng	0.00
4) Naphthalene-d8	9.55	136	105342	5.00	ng	0.00
8) Acenaphthene-d10	13.35	164	52324	5.00	ng	0.00
14) Phenanthrene-d10	16.08	188	145691	5.00	ng	0.00
18) Chrysene-d12	20.21	240	154552	5.00	ng	0.00
24) Perylene-d12	22.10	264	126254	5.00	ng	0.00

System Monitoring Compounds						
2) 2-Fluorophenol	4.93	112	3279	1.30	ng	0.00
3) Phenol-d6	6.56	99	5454	1.30	ng	-0.01
5) Nitrobenzene-d5	8.16	82	5849	0.95	ng	0.00
9) 2,4,6-Tribromophenol	14.89	330	938	0.98	ng	0.00
10) 2-Fluorobiphenyl	12.01	172	11926	1.01	ng	0.00
20) Terphenyl-d14	18.62	244	15582	0.98	ng	0.00

Target Compounds						Qvalue
6) Naphthalene	9.58	128	22605	1.01	ng	99
7) 2-Methylnaphthalene	11.12	142	8907	0.97	ng	94
11) Acenaphthylene	13.10	152	71978	0.97	ng	99
12) Acenaphthene	13.42	154	14862	1.00	ng	84
13) Fluorene	14.40	166	17998	0.98	ng	97
15) Phenanthrene	16.11	178	36127	1.00	ng	100
16) Anthracene	16.20	178	27482	0.99	ng	100
17) Fluoranthene	18.09	202	23846	0.87	ng	100
19) Pyrene	18.48	202	22489	0.94	ng	99
21) Benzo(a)anthracene	20.19	228	19961	0.94	ng	99
22) Chrysene	20.25	228	37812	1.01	ng	98
23) Indeno(1,2,3-cd)pyrene	23.70	276	22591	0.91	ng	95
25) Benzo(b)fluoranthene	21.55	252	23106	0.97	ng	100
26) Benzo(k)fluoranthene	21.58	252	28068	1.00	ng	99
27) Benzo(a)pyrene	22.01	252	20478	0.89	ng	99
28) Dibenzo(a,h)anthracene	23.65	278	17190	0.99	ng	99
29) Benzo(g,h,i)perylene	24.25	276	20884	0.97	ng	99

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

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