

Data Path : Z:\HPCHEM1\BNA E\DATA\BE021916\
 Data File : BE091421.D
 Acq On : 19 Feb 2016 19:32
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001EC

Quant Time: Feb 19 22:39:42 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	6.87	152	41544	5.00	ng	0.00
4) Naphthalene-d8	9.53	136	205321	5.00	ng	-0.01
8) Acenaphthene-d10	13.35	164	133904	5.00	ng	0.00
14) Phenanthrene-d10	16.08	188	330780	5.00	ng	-0.01
18) Chrysene-d12	20.21	240	380908	5.00	ng	0.00
24) Perylene-d12	22.09	264	391953	5.00	ng	0.00

System Monitoring Compounds						
2) 2-Fluorophenol	4.92	112	8932	1.10	ng	-0.03
3) Phenol-d6	6.54	99	13805	1.17	ng	-0.03
5) Nitrobenzene-d5	8.15	82	12907	1.18	ng	-0.02
9) 2,4,6-Tribromophenol	14.89	330	4893	1.05	ng	-0.01
10) 2-Fluorobiphenyl	12.00	172	32977	0.95	ng	-0.01
20) Terphenyl-d14	18.61	244	60545	1.07	ng	0.00

Target Compounds						Qvalue
6) Naphthalene	9.58	128	43338	0.94	ng	100
7) 2-Methylnaphthalene	11.11	142	26282	1.02	ng	98
11) Acenaphthylene	13.09	152	209158	1.04	ng	100
12) Acenaphthene	13.41	154	33589	0.91	ng	97
13) Fluorene	14.40	166	45404	0.96	ng	100
15) Phenanthrene	16.12	178	80642	0.93	ng	99
16) Anthracene	16.20	178	76184	1.05	ng	99
17) Fluoranthene	18.08	202	87318	1.05	ng	99
19) Pyrene	18.47	202	90071	1.09	ng	100
21) Benzo(a)anthracene	20.18	228	83456	1.12	ng	98
22) Chrysene	20.24	228	100452	1.01	ng	98
23) Indeno(1,2,3-cd)pyrene	23.69	276	87825	1.14	ng	100
25) Benzo(b)fluoranthene	21.54	252	84679	0.99	ng	97
26) Benzo(k)fluoranthene	21.57	252	103325	1.07	ng	99
27) Benzo(a)pyrene	22.00	252	87955	1.11	ng	99
28) Dibenzo(a,h)anthracene	23.64	278	82533	1.13	ng	99
29) Benzo(a,h,i)perylene	24.24	276	90918	1.05	ng	98

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

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