

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE021015\
 Data File : BE089567.D
 Acq On : 11 Feb 2015 2:56
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc : MDL-WATER-0.1/0.2
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Feb 11 03:51:26 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIMPAH-BE021015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 11 01:27:13 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.41	152	26594	5.00	ng	0.00
2) Naphthalene-d8	12.32	136	97235	5.00	ng	0.00
6) Acenaphthene-d10	16.50	164	38225	5.00	ng	0.00
11) Phenanthrene-d10	19.83	188	73591	5.00	ng	0.00
16) Chrysene-d12	23.71	240	90798	5.00	ng	0.00
22) Perylene-d12	25.40	264	91981	5.00	ng	0.00

System Monitoring Compounds						
3) Nitrobenzene-d5	10.68	82	6035	0.97	ng	0.00
7) 2-Fluorobiphenyl	14.95	172	10236	0.99	ng	0.00
18) Terphenyl-d14	22.36	244	11351	1.00	ng	0.00

Target Compounds					Qvalue	
4) Naphthalene	12.37	128	21918	1.01	ng	100
5) 2-Methylnaphthalene	14.03	142	12144	1.01	ng	98
8) Acenaphthylene	16.15	152	64218	0.99	ng	100
9) Acenaphthene	16.58	154	10142	0.99	ng	98
10) Fluorene	17.83	166	10974	1.00	ng	98
12) Pentachlorophenol	19.51	266	354	0.92	ng	96
13) Phenanthrene	19.88	178	18008	1.00	ng	99
14) Anthracene	19.97	178	14970	0.98	ng	99
15) Fluoranthene	21.78	202	16967	0.98	ng	99
17) Pyrene	22.10	202	18752	1.03	ng	99
19) Benzo(a)anthracene	23.70	228	67514	0.89	ng	99
20) Chrysene	23.74	228	87397	1.03	ng	100
21) Indeno(1,2,3-cd)pyrene	26.84	276	65040	0.65	ng	99
23) Benzo(b)fluoranthene	24.97	252	9496m	0.73	ng	
24) Benzo(k)fluoranthene	25.00	252	24227m	1.07	ng	
25) Benzo(a)pyrene	25.34	252	11346	0.77	ng	99
26) Dibenzo(a,h)anthracene	26.87	278	12780	0.70	ng	100
27) Benzo(g,h,i)perylene	27.26	276	66227	0.87	ng	97

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

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