

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE010715\
 Data File : BE089350.D
 Acq On : 8 Jan 2015 6:23
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc : LOQ-WATER-1.0PPM-SIM
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Manual Integrations
 APPROVED

TEJASKUMAR
 1/9/2015 5:21:46 PM

Quant Time: Jan 08 15:42:38 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BASE-BE010715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 08 15:21:44 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	12370	5.00	ng	0.00
2) Naphthalene-d8	8.51	136	48428	5.00	ng	0.00
6) Acenaphthene-d10	10.66	164	23043	5.00	ng	0.01
11) Phenanthrene-d10	12.72	188	37139	5.00	ng	0.00
15) Chrysene-d12	18.07	240	22002	5.00	ng	0.00
21) Perylene-d12	21.12	264	18875	5.00	ng	0.00
System Monitoring Compounds						
3) Nitrobenzene-d5	7.65	82	3377	1.00	ng	0.00
7) 2-Fluorobiphenyl	9.85	172	6459	1.03	ng	0.00
17) Terphenyl-d14	15.85	244	4764	1.22	ng	0.00
Target Compounds						Qvalue
4) Naphthalene	8.54	128	11264	1.03	ng	99
5) 2-Methylnaphthalene	9.39	142	7144	1.05	ng	99
8) Acenaphthylene	10.47	152	41315	1.05	ng	99
9) Acenaphthene	10.69	154	6364	1.04	ng	98
10) Fluorene	11.34	166	6949	1.06	ng	100
12) Phenanthrene	12.75	178	39436	1.02	ng	99
13) Anthracene	12.84	178	37770	1.01	ng	99
14) Fluoranthene	14.96	202	8568	1.00	ng	99
16) Pyrene	15.41	202	8874	1.23	ng	99
18) Benzo(a)anthracene	18.05	228	21967	0.98	ng	98
19) Chrysene	18.13	228	22924	1.02	ng	99
20) Indeno(1,2,3-cd)pyrene	23.22	276	19436m	0.91	ng	
22) Benzo(b)fluoranthene	20.37	252	3895	0.94	ng	97
23) Benzo(k)fluoranthene	20.43	252	5354	1.03	ng	98
24) Benzo(a)pyrene	21.01	252	3984	0.98	ng	97
25) Dibenzo(a,h)anthracene	23.31	278	3525	0.95	ng	94
26) Benzo(g,h,i)perylene	23.81	276	18613	0.99	ng	97

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

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