

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE010715\
 Data File : BE089339.D
 Acq On : 7 Jan 2015 21:32
 Operator : TP/IZ
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC0.8

Manual Integrations
 APPROVED

TEJASKUMAR
 1/9/2015 5:21:24 PM

Quant Time: Jan 08 05:44:51 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 05:21:18 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	17382	5.00	ng	0.00
2) Naphthalene-d8	8.51	136	67158	5.00	ng	0.00
6) Acenaphthene-d10	10.66	164	31412	5.00	ng	0.00
11) Phenanthrene-d10	12.72	188	48547	5.00	ng	0.00
15) Chrysene-d12	18.07	240	35810	5.00	ng	0.00
21) Perylene-d12	21.12	264	32891	5.00	ng	0.00
System Monitoring Compounds						
3) Nitrobenzene-d5	7.65	82	3839	0.91	ng	0.00
7) 2-Fluorobiphenyl	9.85	172	7052	0.86	ng	0.00
17) Terphenyl-d14	15.85	244	5196	0.98	ng	0.00
Target Compounds						Qvalue
4) Naphthalene	8.54	128	12609	0.94	ng	99
5) 2-Methylnaphthalene	9.39	142	7936	0.99	ng	96
8) Acenaphthylene	10.47	152	44670	1.00	ng	99
9) Acenaphthene	10.69	154	6925	1.00	ng	90
10) Fluorene	11.34	166	7465	0.99	ng	99
12) Phenanthrene	12.75	178	42246	0.94	ng	99
13) Anthracene	12.84	178	41220	1.00	ng	99
14) Fluoranthene	14.96	202	9348	0.97	ng	100
16) Pyrene	15.40	202	9878	1.13	ng	100
18) Benzo(a)anthracene	18.04	228	29810	1.01	ng	99
19) Chrysene	18.12	228	30540	1.00	ng	99
20) Indeno(1,2,3-cd)pyrene	23.22	276	29029m	0.99	ng	
22) Benzo(b)fluoranthene	20.37	252	5813	0.89	ng	99
23) Benzo(k)fluoranthene	20.42	252	7630	0.91	ng	99
24) Benzo(a)pyrene	21.01	252	5775	0.93	ng	100
25) Dibenzo(a,h)anthracene	23.30	278	5378	0.83	ng	100
26) Benzo(g,h,i)perylene	23.80	276	27096	0.92	ng	100

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

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