

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE010715\
 Data File : BE089341.D
 Acq On : 7 Jan 2015 22:38
 Operator : TP/IZ
 Sample : SSTDICC002
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC002

Manual Integrations
 APPROVED

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

Quant Time: Jan 08 05:48:00 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	16824	5.00	ng	0.00
2) Naphthalene-d8	8.51	136	64394	5.00	ng	0.00
6) Acenaphthene-d10	10.65	164	29451	5.00	ng	0.00
11) Phenanthrene-d10	12.72	188	45771	5.00	ng	0.00
15) Chrysene-d12	18.07	240	35497	5.00	ng	0.00
21) Perylene-d12	21.12	264	32467	5.00	ng	0.00
System Monitoring Compounds						
3) Nitrobenzene-d5	7.65	82	9578	2.36	ng	0.00
7) 2-Fluorobiphenyl	9.85	172	16485	2.15	ng	0.00
17) Terphenyl-d14	15.85	244	12563	2.40	ng	0.00
Target Compounds						Qvalue
4) Naphthalene	8.53	128	28679	2.22	ng	99
5) 2-Methylnaphthalene	9.39	142	17888	2.32	ng	95
8) Acenaphthylene	10.47	152	101186	2.42	ng	98
9) Acenaphthene	10.69	154	15216	2.35	ng	94
10) Fluorene	11.34	166	16903	2.39	ng	99
12) Phenanthrene	12.75	178	95149	2.25	ng	99
13) Anthracene	12.84	178	94734	2.44	ng	98
14) Fluoranthene	14.96	202	21657	2.37	ng	98
16) Pyrene	15.40	202	22995	2.65	ng	99
18) Benzo(a)anthracene	18.05	228	70826	2.42	ng	99
19) Chrysene	18.12	228	71867	2.37	ng	99
20) Indeno(1,2,3-cd)pyrene	23.21	276	74142m	2.54	ng	
22) Benzo(b)fluoranthene	20.36	252	14818	2.30	ng	99
23) Benzo(k)fluoranthene	20.42	252	17940	2.16	ng	99
24) Benzo(a)pyrene	21.00	252	14427	2.36	ng	98
25) Dibenzo(a,h)anthracene	23.30	278	13344	2.07	ng	97
26) Benzo(g,h,i)perylene	23.79	276	65818	2.26	ng	99

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

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