

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE010715\
 Data File : BE089342.D
 Acq On : 7 Jan 2015 23:11
 Operator : TP/IZ
 Sample : SSTDICC005
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

TEJASKUMAR
 1/9/2015 5:21:18 PM

Quant Time: Jan 08 05:48:39 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	17616	5.00	ng	0.00
2) Naphthalene-d8	8.51	136	67455	5.00	ng	0.00
6) Acenaphthene-d10	10.65	164	31224	5.00	ng	0.00
11) Phenanthrene-d10	12.72	188	49294	5.00	ng	0.00
15) Chrysene-d12	18.07	240	36087	5.00	ng	0.00
21) Perylene-d12	21.12	264	31984	5.00	ng	0.00
System Monitoring Compounds						
3) Nitrobenzene-d5	7.65	82	25033	5.89	ng	0.00
7) 2-Fluorobiphenyl	9.85	172	41194	5.07	ng	0.00
17) Terphenyl-d14	15.84	244	31423	5.90	ng	0.00
Target Compounds						Qvalue
4) Naphthalene	8.53	128	72353	5.35	ng	100
5) 2-Methylnaphthalene	9.39	142	45652	5.64	ng	98
8) Acenaphthylene	10.47	152	260636	5.89	ng	99
9) Acenaphthene	10.69	154	39016	5.68	ng	99
10) Fluorene	11.34	166	43978	5.86	ng	99
12) Phenanthrene	12.75	178	244433	5.36	ng	99
13) Anthracene	12.84	178	245069	5.87	ng	100
14) Fluoranthene	14.95	202	55274	5.63	ng	99
16) Pyrene	15.40	202	58358	6.61	ng	100
18) Benzo(a)anthracene	18.05	228	178870	6.01	ng	99
19) Chrysene	18.12	228	176539	5.72	ng	99
20) Indeno(1,2,3-cd)pyrene	23.21	276	179491m	6.05	ng	
22) Benzo(b)fluoranthene	20.36	252	36926	5.81	ng	98
23) Benzo(k)fluoranthene	20.42	252	44221	5.40	ng	98
24) Benzo(a)pyrene	21.00	252	36471	6.06	ng	97
25) Dibenzo(a,h)anthracene	23.29	278	33958	5.36	ng	97
26) Benzo(g,h,i)perylene	23.79	276	163200	5.70	ng	98

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

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