

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE032415\
 Data File : BE089817.D
 Acq On : 24 Mar 2015 16:10
 Operator : TP/JJ
 Sample : SSTDICC005
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

mohammad
 3/25/2015 5:43:28 PM

Quant Time: Mar 24 16:54:49 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIMPAH-BE032415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 24 15:59:18 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	175256	5.00	ng	0.00
4) Naphthalene-d8	10.26	136	726077	5.00	ng	0.00
8) Acenaphthene-d10	14.40	164	330849	5.00	ng	0.01
13) Phenanthrene-d10	17.92	188	566052	5.00	ng	0.00
19) Chrysene-d12	24.26	240	555393	5.00	ng	0.00
25) Perylene-d12	27.41	264	494689	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.68	82	254361	6.76	ng	0.00
9) 2-Fluorobiphenyl	12.85	172	406515	5.34	ng	0.00
21) Terphenyl-d14	21.86	244	399756	4.85	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.54	88	108913	4.39	ng	90
3) n-Nitrosodimethylamine	2.95	42	186909	5.85	ng	# 89
6) Naphthalene	10.31	128	692874	4.77	ng	96
7) 2-Methylnaphthalene	11.93	142	429751	5.53	ng	97
10) Acenaphthylene	14.04	152	2533349	5.49	ng	99
11) Acenaphthene	14.46	154	364195	5.04	ng	85
12) Fluorene	15.71	166	449898	5.17	ng	94
14) Hexachlorobenzene	16.98	284	92914	4.92	ng	97
15) Pentachlorophenol	17.50	266	28725	0.27	ng	94
16) Phenanthrene	17.97	178	618496	4.97	ng	100
17) Anthracene	18.10	178	602404	5.59	ng	99
18) Fluoranthene	20.82	202	642520	6.07	ng	99
20) Pyrene	21.34	202	666212	4.64	ng	98
22) Benzo(a)anthracene	24.24	228	571264	5.73	ng	98
23) Chrysene	24.32	228	583821	5.21	ng	99
24) Indeno(1,2,3-cd)pyrene	29.38	276	527267m	10.94	ng	
26) Benzo(b)fluoranthene	26.64	252	495424	7.80	ng	# 98
27) Benzo(k)fluoranthene	26.70	252	602253m	6.34	ng	
28) Benzo(a)pyrene	27.30	252	470707	7.86	ng	# 89
29) Dibenzo(a,h)anthracene	29.44	278	421705	11.31	ng	# 77
30) Benzo(g,h,i)perylene	29.83	276	501536	7.68	ng	# 88

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

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