

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020315\
 Data File : BE089465.D
 Acq On : 3 Feb 2015 15:40
 Operator : TP/IZ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

apatel
 2/4/2015 2:54:46 PM

Quant Time: Feb 04 00:34:56 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE020315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 04 00:22:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.53	152	59207	5.00	ng	0.00
7) Naphthalene-d8	12.43	136	234127	5.00	ng	0.00
13) Acenaphthene-d10	16.62	164	100469	5.00	ng	0.00
17) Phenanthrene-d10	19.94	188	193248	5.00	ng	0.00
20) Chrysene-d12	23.79	240	189871	5.00	ng	0.00
27) Perylene-d12	25.49	264	180237	5.00	ng	0.00
System Monitoring Compounds						
3) 2-Fluorophenol	6.72	112	193553	9.77	ng	0.00
4) Phenol-d6	8.85	99	239535	10.30	ng	0.00
8) Nitrobenzene-d5	10.79	82	187277	10.96	ng	0.00
14) 2,4,6-Tribromophenol	18.52	330	44014	16.76	ng	0.01
15) 2-Fluorobiphenyl	15.07	172	262662	9.37	ng	0.00
22) Terphenyl-d14	22.43	244	283211	12.94	ng	0.00
Target Compounds						Qvalue
2) n-Nitrosodimethylamine	3.52	42	121108	12.72	ng	# 79
5) Phenol	8.87	94	257593m	10.31	ng	
6) bis(2-Chloroethyl)ether	9.01	93	184343	10.11	ng	# 99
9) Nitrobenzene	10.84	77	211565	10.91	ng	98
10) 2,4-Dimethylphenol	11.78	122	152747	10.98	ng	98
11) 2,4-Dichlorophenol	12.16	162	115371	11.00	ng	90
12) Hexachlorobutadiene	12.81	225	69576	9.43	ng	99
16) 2,4-Dinitrophenol	16.85	184	18482	30.57	ng	# 78
18) Hexachlorobenzene	19.18	284	96384	9.50	ng	98
19) Pentachlorophenol	19.61	266	37402	31.38	ng	97
21) Benzidine	22.10	184	94872	7.01	ng	99
23) Benzo(a)anthracene	23.78	228	1775482	11.59	ng	97
24) 3,3'-Dichlorobenzidine	23.77	252	160472	11.91	ng	# 86
25) Chrysene	23.82	228	1665393	9.07	ng	93
26) Indeno(1,2,3-cd)pyrene	26.97	276	498940	9.28	ng	99
28) Benzo(b)fluoranthene	25.05	252	413295	16.80	ng	98
29) Benzo(k)fluoranthene	25.08	252	459952	10.17	ng	99
30) Benzo(a)pyrene	25.42	252	407183	14.76	ng	97
31) Dibenzo(a,h)anthracene	27.00	278	428190	12.95	ng	98

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

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