

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020315\
 Data File : BE089467.D
 Acq On : 3 Feb 2015 16:58
 Operator : TP/IZ
 Sample : SSTDICC002
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC002

Manual Integrations
 APPROVED

apatel
 2/4/2015 2:54:53 PM

Quant Time: Feb 04 00:37:50 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	55148	5.00	ng	0.00
7) Naphthalene-d8	12.43	136	218731	5.00	ng	0.00
13) Acenaphthene-d10	16.61	164	94634	5.00	ng	0.00
17) Phenanthrene-d10	19.93	188	180609	5.00	ng	0.00
20) Chrysene-d12	23.78	240	181480	5.00	ng	0.00
27) Perylene-d12	25.49	264	170874	5.00	ng	0.00
System Monitoring Compounds						
3) 2-Fluorophenol	6.72	112	36702	1.99	ng	0.00
4) Phenol-d6	8.84	99	46044	2.13	ng	0.00
8) Nitrobenzene-d5	10.79	82	34642	2.17	ng	0.00
14) 2,4,6-Tribromophenol	18.51	330	7276	2.94	ng	0.00
15) 2-Fluorobiphenyl	15.06	172	51339	1.94	ng	0.00
22) Terphenyl-d14	22.43	244	54401	2.60	ng	0.00
Target Compounds						Qvalue
2) n-Nitrosodimethylamine	3.54	42	21812	2.46	ng	# 96
5) Phenol	8.86	94	50184m	2.16	ng	
6) bis(2-Chloroethyl)ether	9.00	93	35976	2.12	ng	# 100
9) Nitrobenzene	10.83	77	40125	2.22	ng	100
10) 2,4-Dimethylphenol	11.77	122	28572	2.20	ng	95
11) 2,4-Dichlorophenol	12.15	162	21169	2.16	ng	97
12) Hexachlorobutadiene	12.81	225	13860	2.01	ng	99
16) 2,4-Dinitrophenol	16.84	184	1209	2.12	ng	86
18) Hexachlorobenzene	19.17	284	19359	2.04	ng	94
19) Pentachlorophenol	19.60	266	3328	2.99	ng	99
21) Benzidine	22.09	184	18753	1.45	ng	98
23) Benzo(a)anthracene	23.77	228	334724	2.29	ng	99
24) 3,3'-Dichlorobenzidine	23.76	252	29442	2.29	ng	99
25) Chrysene	23.82	228	351077	2.00	ng	99
26) Indeno(1,2,3-cd)pyrene	26.97	276	91593	1.78	ng	99
28) Benzo(b)fluoranthene	25.04	252	70079	3.01	ng	99
29) Benzo(k)fluoranthene	25.07	252	98655	2.30	ng	99
30) Benzo(a)pyrene	25.42	252	70624	2.70	ng	98
31) Dibenzo(a,h)anthracene	26.99	278	75914	2.42	ng	99

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

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