

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020315\
 Data File : BE089482.D
 Acq On : 4 Feb 2015 3:25
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Manual Integrations
 APPROVED

apatel
 2/4/2015 2:55:34 PM

Quant Time: Feb 04 10:51:24 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 00:53:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.53	152	51572	5.00	ng	0.00
7) Naphthalene-d8	12.43	136	210027	5.00	ng	0.00
13) Acenaphthene-d10	16.61	164	91448	5.00	ng	0.00
17) Phenanthrene-d10	19.93	188	171762	5.00	ng	0.00
20) Chrysene-d12	23.78	240	164308	5.00	ng	0.00
27) Perylene-d12	25.49	264	149532	5.00	ng	0.00
System Monitoring Compounds						
3) 2-Fluorophenol	6.73	112	15775	0.95	ng	0.00
4) Phenol-d6	8.85	99	20380	0.97	ng	0.00
8) Nitrobenzene-d5	10.79	82	15127	0.95	ng	0.00
14) 2,4,6-Tribromophenol	18.52	330	2635	0.83	ng	0.00
15) 2-Fluorobiphenyl	15.06	172	23899	0.99	ng	0.00
22) Terphenyl-d14	22.43	244	23759	1.02	ng	0.00
Target Compounds						Qvalue
2) n-Nitrosodimethylamine	3.56	42	9280	0.96	ng	# 70
5) Phenol	8.87	94	22134m	0.97	ng	
6) bis(2-Chloroethyl)ether	9.01	93	16052	0.97	ng	# 100
9) Nitrobenzene	10.83	77	17730	0.98	ng	99
10) 2,4-Dimethylphenol	11.78	122	12534	0.95	ng	92
11) 2,4-Dichlorophenol	12.16	162	9262	0.95	ng	89
12) Hexachlorobutadiene	12.81	225	6455	0.99	ng	99
16) 2,4-Dinitrophenol	16.85	184	311m	0.92	ng	
18) Hexachlorobenzene	19.17	284	8862	1.00	ng	99
19) Pentachlorophenol	19.61	266	858m	0.81	ng	
21) Benzidine	22.10	184	16967	2.15	ng	99
23) Benzo(a)anthracene	23.77	228	143943	0.99	ng	98
24) 3,3'-Dichlorobenzidine	23.77	252	11423	0.94	ng	# 90
25) Chrysene	23.82	228	151373	1.00	ng	98
26) Indeno(1,2,3-cd)pyrene	26.97	276	34059	0.91	ng	98
28) Benzo(b)fluoranthene	25.04	252	23134	0.84	ng	99
29) Benzo(k)fluoranthene	25.07	252	42898	1.11	ng	99
30) Benzo(a)pyrene	25.42	252	25188	0.91	ng	99
31) Dibenzo(a,h)anthracene	27.00	278	26415	0.88	ng	99

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

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