

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE012915\
 Data File : BE089411.D
 Acq On : 29 Jan 2015 12:02
 Operator : TP/IZ
 Sample : SSTDICC0.5
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC0.5

Manual Integrations
 APPROVED

apatel
 1/30/2015 5:55:26 PM

Quant Time: Jan 30 02:51:56 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE012915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 30 02:34:28 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.58	152	142382	5.00	ng	0.00
7) Naphthalene-d8	12.50	136	537656	5.00	ng	0.00
13) Acenaphthene-d10	16.68	164	211174	5.00	ng	0.00
17) Phenanthrene-d10	19.99	188	420156	5.00	ng	0.00
20) Chrysene-d12	23.84	240	516859	5.00	ng	0.00
27) Perylene-d12	25.55	264	523057	5.00	ng	0.00
System Monitoring Compounds						
3) 2-Fluorophenol	6.74	112	25757	0.56	ng	0.00
4) Phenol-d6	8.85	99	31480	0.56	ng	0.00
8) Nitrobenzene-d5	10.84	82	21000	0.54	ng	0.00
14) 2,4,6-Tribromophenol	18.58	330	2660	0.54	ng	0.00
15) 2-Fluorobiphenyl	15.13	172	29006	0.46	ng	0.00
22) Terphenyl-d14	22.48	244	30012	0.53	ng	0.00
Target Compounds						Qvalue
2) n-Nitrosodimethylamine	3.60	42	13927	0.57	ng	# 86
5) Phenol	8.87	94	34673	0.58	ng	94
6) bis(2-Chloroethyl)ether	9.06	93	25745	0.58	ng	# 99
9) Nitrobenzene	10.89	77	24296	0.56	ng	99
10) 2,4-Dimethylphenol	11.82	122	17189	0.52	ng	97
11) 2,4-Dichlorophenol	12.21	162	11660	0.46	ng	97
12) Hexachlorobutadiene	12.89	225	8037	0.45	ng	99
16) 2,4-Dinitrophenol	16.90	184	201	0.14	ng	# 70
18) Hexachlorobenzene	19.23	284	10150	0.52	ng	95
19) Pentachlorophenol	19.65	266	1101	0.29	ng	97
21) Benzidine	22.14	184	13445	0.32	ng	97
23) Benzo(a)anthracene	23.83	228	180998	0.39	ng	94
24) 3,3'-Dichlorobenzidine	23.81	252	12102	0.32	ng	# 96
25) Chrysene	23.87	228	266794	0.46	ng	100
26) Indeno(1,2,3-cd)pyrene	27.05	276	28652	0.21	ng	# 100
28) Benzo(b)fluoranthene	25.09	252	16595m	0.25	ng	
29) Benzo(k)fluoranthene	25.13	252	65841m	0.44	ng	
30) Benzo(a)pyrene	25.48	252	21851	0.27	ng	98
31) Dibenzo(a,h)anthracene	27.08	278	20005	0.24	ng	97

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

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