

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE010515\
 Data File : BE089324.D
 Acq On : 5 Jan 2015 20:57
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Manual Integrations
 APPROVED

TEJASKUMAR
 1/7/2015 9:23:51 AM

Quant Time: Jan 06 14:12:11 2015
 Quant Method : Z:\HPCHEM1\BNA_E\methods\SIM-BE121914.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 20 00:43:57 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	30382	5.00	ng	0.00
7) Naphthalene-d8	8.51	136	116267	5.00	ng	0.00
13) Acenaphthene-d10	10.66	164	52925	5.00	ng	0.00
17) Phenanthrene-d10	12.71	188	77503	5.00	ng	0.00
20) Chrysene-d12	18.06	240	55052	5.00	ng	0.00
27) Perylene-d12	21.11	264	50181	5.00	ng	0.00
System Monitoring Compounds						
3) 2-Fluorophenol	5.32	112	8135	1.23	ng	0.05
4) Phenol-d6	6.62	99	10729	1.08	ng	0.07
8) Nitrobenzene-d5	7.64	82	9001	1.03	ng	0.00
14) 2,4,6-Tribromophenol	11.68	330	1599	1.12	ng	0.00
15) 2-Fluorobiphenyl	9.85	172	15317	1.02	ng	0.00
22) Terphenyl-d14	15.84	244	9877	1.00	ng	0.00
Target Compounds						Qvalue
2) n-Nitrosodimethylamine	2.69	42	4908	1.24	ng	# 89
5) Phenol	6.64	94	13465m	1.13	ng	
6) bis(2-Chloroethyl)ether	6.65	93	10510	1.15	ng	# 92
9) Nitrobenzene	7.67	77	10392	1.15	ng	99
10) 2,4-Dimethylphenol	8.18	122	8109	1.20	ng	# 80
11) 2,4-Dichlorophenol	8.40	162	6790	1.20	ng	93
12) Hexachlorobutadiene	8.69	225	3482	1.09	ng	99
16) 2,4-Dinitrophenol	10.79	184	430	0.75	ng	# 75
18) Hexachlorobenzene	12.10	284	4160	1.12	ng	97
19) Pentachlorophenol	12.45	266	1228	1.39	ng	94
21) Benzidine	15.34	184	6462	0.93	ng	97
23) Benzo(a)anthracene	18.04	228	58339	1.19	ng	99
24) 3,3'-Dichlorobenzidine	18.11	252	5288	1.34	ng	98
25) Chrysene	18.12	228	57920	1.12	ng	98
26) Indeno(1,2,3-cd)pyrene	23.20	276	55069m	1.37	ng	
28) Benzo(b)fluoranthene	20.36	252	12459	1.33	ng	100
29) Benzo(k)fluoranthene	20.42	252	14275	1.12	ng	100
30) Benzo(a)pyrene	21.00	252	12330	1.34	ng	99
31) Dibenzo(a,h)anthracene	23.29	278	10906	1.39	ng	97

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

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