

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE012815\
 Data File : BE089399.D
 Acq On : 28 Jan 2015 21:48
 Operator : TP/IZ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC010

Quant Time: Jan 29 12:13:41 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE012815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 29 12:06:13 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.58	152	84323	5.00	ng	0.00
7) Naphthalene-d8	12.50	136	254976	5.00	ng	0.00
13) Acenaphthene-d10	16.68	164	75016	5.00	ng	0.00
17) Phenanthrene-d10	19.98	188	103632	5.00	ng	0.00
20) Chrysene-d12	23.84	240	227100	5.00	ng	0.00
27) Perylene-d12	25.55	264	355398	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 2-Fluorophenol	6.72	112	256686	9.74	ng	0.00
4) Phenol-d6	8.84	99	289788	8.92	ng	0.00
8) Nitrobenzene-d5	10.84	82	182634	10.16	ng	0.00
14) 2,4,6-Tribromophenol	18.58	330	14736	0.23	ng	0.00
15) 2-Fluorobiphenyl	15.13	172	243574	0.31	ng	0.00
22) Terphenyl-d14	22.48	244	170892	6.79	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) n-Nitrosodimethylamine	3.54	42	147195	9.77	ng	# 85
5) Phenol	8.87	94	307926	8.92	ng	# 51
6) bis(2-Chloroethyl)ether	9.06	93	228278	8.89	ng	# 99
9) Nitrobenzene	10.89	77	206838	10.19	ng	99
10) 2,4-Dimethylphenol	11.82	122	149878	9.69	ng	99
11) 2,4-Dichlorophenol	12.20	162	116738	9.81	ng	88
12) Hexachlorobutadiene	12.89	225	83905	9.93	ng	99
16) 2,4-Dinitrophenol	16.89	184	8569	2.19	ng	92
18) Hexachlorobenzene	19.23	284	49762	10.43	ng	99
19) Pentachlorophenol	19.65	266	17350	18.83	ng	99
21) Benzidine	22.14	184	233176	8.70	ng	98
23) Benzo(a)anthracene	23.83	228	2106827	10.02	ng	98
24) 3,3'-Dichlorobenzidine	23.81	252	251513	15.02	ng	98
25) Chrysene	23.87	228	2410374	9.37	ng	97
26) Indeno(1,2,3-cd)pyrene	27.07	276	1245986	20.78	ng	# 100
28) Benzo(b)fluoranthene	25.10	252	635514	13.61	ng	99
29) Benzo(k)fluoranthene	25.13	252	1042211	9.21	ng	98
30) Benzo(a)pyrene	25.48	252	879876	16.05	ng	97
31) Dibenzo(a,h)anthracene	27.10	278	966660	16.58	ng	# 93

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

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