

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE012915\
 Data File : BE089423.D
 Acq On : 29 Jan 2015 21:02
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Jan 30 07:27:24 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 03:48:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.58	152	165772	5.00	ng	0.00
7) Naphthalene-d8	12.50	136	616047	5.00	ng	0.00
13) Acenaphthene-d10	16.68	164	245613	5.00	ng	0.00
17) Phenanthrene-d10	19.99	188	504580	5.00	ng	0.00
20) Chrysene-d12	23.84	240	665818	5.00	ng	0.00
27) Perylene-d12	25.55	264	706291	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 2-Fluorophenol	6.74	112	57817	0.93	ng	0.00
4) Phenol-d6	8.85	99	69559	0.94	ng	0.00
8) Nitrobenzene-d5	10.85	82	46750	0.93	ng	0.00
14) 2,4,6-Tribromophenol	18.58	330	6195	0.91	ng	0.00
15) 2-Fluorobiphenyl	15.13	172	64633	0.94	ng	0.00
22) Terphenyl-d14	22.48	244	82902	1.08	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) n-Nitrosodimethylamine	3.58	42	33004	1.02	ng	83
5) Phenol	8.88	94	75746	0.93	ng	93
6) bis(2-Chloroethyl)ether	9.06	93	55868	0.93	ng	# 99
9) Nitrobenzene	10.89	77	54409	0.95	ng	99
10) 2,4-Dimethylphenol	11.83	122	38505	0.95	ng	94
11) 2,4-Dichlorophenol	12.21	162	26911	0.95	ng	90
12) Hexachlorobutadiene	12.89	225	17711	0.93	ng	99
16) 2,4-Dinitrophenol	16.90	184	1451	0.81	ng	87
18) Hexachlorobenzene	19.23	284	23518	0.95	ng	94
19) Pentachlorophenol	19.65	266	3158	0.68	ng	95
21) Benzidine	22.14	184	46650	0.80	ng	99
23) Benzo(a)anthracene	23.83	228	558596	1.10	ng	100
24) 3,3'-Dichlorobenzidine	23.81	252	46615	1.17	ng	97
25) Chrysene	23.87	228	639211	0.92	ng	97
26) Indeno(1,2,3-cd)pyrene	27.07	276	152142	1.27	ng	# 100
28) Benzo(b)fluoranthene	25.10	252	95457	1.37	ng	98
29) Benzo(k)fluoranthene	25.13	252	197621	1.07	ng	99
30) Benzo(a)pyrene	25.48	252	106620	1.24	ng	99
31) Dibenzo(a,h)anthracene	27.09	278	119561	1.39	ng	96

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

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