

Data Path : Z:\HPCHEM1\BNA E\DATA\BE021615\
 Data File : BE089581.D
 Acq On : 16 Feb 2015 23:03
 Operator : TP/IZ
 Sample : PB81705BSD
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB81705BSD

Quant Time: Feb 17 00:23:50 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SIM-BE021615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 16 18:23:16 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.34	152	81930	5.00	ng	0.00
7) Naphthalene-d8	12.25	136	313500	5.00	ng	0.00
13) Acenaphthene-d10	16.42	164	151417	5.00	ng	0.00
17) Phenanthrene-d10	19.78	188	293118	5.00	ng	0.00
20) Chrysene-d12	23.66	240	390314	5.00	ng	0.00
27) Perylene-d12	25.34	264	357370	5.00	ng	0.00

System Monitoring Compounds						
3) 2-Fluorophenol	6.51	112	259287	10.95	ng	0.00
4) Phenol-d6	8.63	99	345229	10.99	ng	0.00
8) Nitrobenzene-d5	10.61	82	154588	11.60	ng	0.00
14) 2,4,6-Tribromophenol	18.32	330	55945	16.40	ng	0.00
15) 2-Fluorobiphenyl	14.88	172	364851	8.92	ng	0.00
22) Terphenyl-d14	22.31	244	487435	9.06	ng	0.00

Target Compounds						Qvalue
2) n-Nitrosodimethylamine	3.31	42	90940	7.00	ng	# 59
5) Phenol	8.66	94	257364	7.62	ng	# 79
6) bis(2-Chloroethyl)ether	8.83	93	211913	8.24	ng	# 92
9) Nitrobenzene	10.65	77	180612	11.22	ng	94
10) 2,4-Dimethylphenol	11.59	122	143591	7.77	ng	100
11) 2,4-Dichlorophenol	11.97	162	130411	9.02	ng	99
12) Hexachlorobutadiene	12.64	225	78768	8.27	ng	100
16) 2,4-Dinitrophenol	16.66	184	29527	56.07	ng	# 62
18) Hexachlorobenzene	19.00	284	109800	8.91	ng	92
19) Pentachlorophenol	19.44	266	82545	36.91	ng	96
21) Benzidine	21.97	184	788717	22.57	ng	99
23) Benzo(a)anthracene	23.64	228	3506007	9.07	ng	93
24) 3,3'-Dichlorobenzidine	23.63	252	316067	11.26	ng	96
25) Chrysene	23.69	228	3279131	8.83	ng	95
26) Indeno(1,2,3-cd)pyrene	26.76	276	955979	9.91	ng	99
28) Benzo(b)fluoranthene	24.92	252	814057	10.52	ng	97
29) Benzo(k)fluoranthene	24.95	252	916163	9.08	ng	99
30) Benzo(a)pyrene	25.28	252	792692	10.69	ng	99
31) Dibenzo(a,h)anthracene	26.78	278	797830	10.56	ng	99

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

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