

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE021715\
 Data File : BE089605.D
 Acq On : 17 Feb 2015 22:20
 Operator : TP/IZ
 Sample : PB81705BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB81705BS

Quant Time: Feb 18 14:43:20 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE021715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 18 14:11:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.34	152	63216	5.00	ng	0.00
7) Naphthalene-d8	12.24	136	253699	5.00	ng	0.00
13) Acenaphthene-d10	16.42	164	132450	5.00	ng	0.00
17) Phenanthrene-d10	19.77	188	262292	5.00	ng	0.00
20) Chrysene-d12	23.66	240	303433	5.00	ng	0.00
27) Perylene-d12	25.34	264	252752	5.00	ng	0.00

System Monitoring Compounds						
3) 2-Fluorophenol	6.51	112	220561	13.40	ng	0.00
4) Phenol-d6	8.63	99	295948	13.66	ng	0.00
8) Nitrobenzene-d5	10.61	82	142848	10.16	ng	0.00
14) 2,4,6-Tribromophenol	18.32	330	55606	14.29	ng	0.00
15) 2-Fluorobiphenyl	14.88	172	347244	9.81	ng	0.00
22) Terphenyl-d14	22.30	244	458295	11.19	ng	0.00

Target Compounds						Qvalue
2) n-Nitrosodimethylamine	3.32	42	66087	7.91	ng	# 61
5) Phenol	8.66	94	203064	8.70	ng	84
6) bis(2-Chloroethyl)ether	8.83	93	165146	8.88	ng	# 98
9) Nitrobenzene	10.65	77	150976	8.93	ng	96
10) 2,4-Dimethylphenol	11.58	122	117283	8.98	ng	94
11) 2,4-Dichlorophenol	11.96	162	108255	10.31	ng	97
12) Hexachlorobutadiene	12.63	225	63289	8.07	ng	100
16) 2,4-Dinitrophenol	16.66	184	24457	31.26	ng	# 62
18) Hexachlorobenzene	19.00	284	99910	8.83	ng	# 87
19) Pentachlorophenol	19.44	266	74732	13.73	ng	98
21) Benzidine	21.96	184	458981	20.64	ng	100
23) Benzo(a)anthracene	23.64	228	2723658	10.53	ng	98
24) 3,3'-Dichlorobenzidine	23.64	252	233219	8.36	ng	# 97
25) Chrysene	23.69	228	2602251	8.79	ng	95
26) Indeno(1,2,3-cd)pyrene	26.75	276	630604	9.03	ng	# 99
28) Benzo(b)fluoranthene	24.92	252	572287	9.96	ng	99
29) Benzo(k)fluoranthene	24.95	252	691194	8.66	ng	98
30) Benzo(a)pyrene	25.28	252	560618	10.16	ng	99
31) Dibenzo(a,h)anthracene	26.78	278	526628	9.76	ng	100

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

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