

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062615\
 Data File : BE089948.D
 Acq On : 26 Jun 2015 18:23
 Operator : TP/IZ
 Sample : G2782-31MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X1SS0610204-150624MS

Quant Time: Jun 27 01:00:16 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 25 16:51:27 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	26289	5.00	ng	0.00
6) Naphthalene-d8	10.39	136	119908	5.00	ng	0.00
12) Acenaphthene-d10	14.24	164	70208	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	162823	5.00	ng	0.00
25) Chrysene-d12	21.13	240	196925	5.00	ng	0.00
32) Perylene-d12	23.46	264	177036	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.26	112	51833	10.92	ng	0.00
5) Phenol-d6	6.81	99	85527	11.76	ng	0.00
7) Nitrobenzene-d5	8.77	82	65777	7.37	ng	0.00
13) 2,4,6-Tribromophenol	15.74	330	13670	10.92	ng	0.00
14) 2-Fluorobiphenyl	12.87	172	147641	7.22	ng	0.00
27) Terphenyl-d14	19.60	244	227525	6.69	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.26	88	10164	3.23	ng	99
3) n-Nitrosodimethylamine	3.55	42	15496	3.28	ng	99
8) Nitrobenzene	8.81	77	34757	3.66	ng	97
9) Naphthalene	10.44	128	90312	3.39	ng	99
10) Hexachlorobutadiene	10.74	225	14118	3.35	ng	99
11) 2-Methylnaphthalene	12.05	142	62836	3.44	ng	99
15) Acenaphthylene	13.96	152	429448	3.49	ng	100
16) Acenaphthene	14.30	154	62942	3.48	ng	98
17) Fluorene	15.29	166	85786	3.79	ng	98
19) 4-Bromophenyl-phenylether	16.18	248	23992	3.61	ng	94
20) Hexachlorobenzene	16.30	284	104640	3.59	ng	98
21) Pentachlorophenol	16.63	266	8807	5.65	ng	90
22) Phenanthrene	17.01	178	141370	3.46	ng	100
23) Anthracene	17.12	178	140937	3.72	ng	99
24) Fluoranthene	19.02	202	164835	3.72	ng	100
26) Pyrene	19.38	202	177956	3.57	ng	99
28) Benzo(a)anthracene	21.12	228	174446	3.71	ng	98
29) Chrysene	21.17	228	179172	3.48	ng	98
30) Bis(2-ethylhexyl)phthalate	21.07	149	25256	6.30	ng	100
31) Indeno(1,2,3-cd)pyrene	25.88	276	159812	4.23	ng	# 100
33) Benzo(b)fluoranthene	22.75	252	149338	3.99	ng	99
34) Benzo(k)fluoranthene	22.80	252	164163	3.87	ng	99
35) Benzo(a)pyrene	23.36	252	141026	4.24	ng	99
36) Dibenzo(a,h)anthracene	25.90	278	130079	4.20	ng	99
37) Benzo(g,h,i)perylene	26.62	276	136135	4.09	ng	99

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

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