

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE050719\
 Data File : BE099570.D
 Acq On : 7 May 2019 17:59
 Operator : JU/SJ
 Sample : K2707-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 BP-TT-MW201D1-GW-20190502

Quant Time: May 08 01:37:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE050719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 07 14:10:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	7	0.40	ng	-0.01
7) Naphthalene-d8	10.63	136	28	0.40	ng	0.00
13) Acenaphthene-d10	14.51	164	26	0.40	ng	0.01
19) Phenanthrene-d10	17.26	188	8	0.40	ng	0.01
27) Chrysene-d12	21.41	240	81	0.40	ng	-0.01
34) Perylene-d12	23.96	264	3	0.40	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.35	112	17	0.88	ng	-0.04
5) Phenol-d6	7.03	99	22	0.86	ng	0.05
8) Nitrobenzene-d5	8.98	82	19	0.70	ng	0.00
11) 2-Methylnaphthalene-d10	12.19	152	7	0.14	ng	-0.03
14) 2,4,6-Tribromophenol	15.97	330	8	0.46	ng	-0.01
15) 2-Fluorobiphenyl	13.12	172	16	0.16	ng	0.00
25) Fluoranthene-d10	19.28	212	41	0.38	ng	0.00
29) Terphenyl-d14	19.86	244	7	0.05	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.32	88	7	0.708	ng	# 1
3) n-Nitrosodimethylamine	3.57	42	55	8.405	ng	# 1
6) bis(2-Chloroethyl)ether	7.25	93	12	0.564	ng	# 1
9) Naphthalene	10.69	128	53	0.179	ng	# 46
10) Hexachlorobutadiene	10.93	225	4	0.186	ng	# 17
12) 2-Methylnaphthalene	12.34	142	13	0.259	ng	# 42
16) Acenaphthylene	14.17	152	16	0.128	ng	# 86
18) Fluorene	15.57	166	10	0.100	ng	# 9
20) 4-Bromophenyl-phenylether	16.45	248	14	3.080	ng	# 65
21) Hexachlorobenzene	16.59	284	4	0.885	ng	# 46
22) Pentachlorophenol	16.88	266	6	3.225	ng	# 67
23) Phenanthrene	17.32	178	35	1.775	ng	# 56
24) Anthracene	17.42	178	22	1.160	ng	# 60
26) Fluoranthene	19.31	202	25	0.823	ng	# 88
28) Pyrene	19.67	202	23	0.106	ng	# 78
30) Benzo(a)anthracene	21.43	228	10	0.040	ng	# 1
31) Chrysene	21.46	228	22	0.089	ng	# 1
32) Bis(2-ethylhexyl)phthalate	21.33	149	128	0.322	ng	# 54
33) Indeno(1,2,3-cd)pyrene	26.67	276	6	0.020	ng	# 57
35) Benzo(b)fluoranthene	23.16	252	44	5.317	ng	# 1
36) Benzo(k)fluoranthene	23.21	252	26	3.193	ng	# 1
37) Benzo(a)pyrene	23.84	252	5	0.627	ng	# 1
38) Dibenzo(a,h)anthracene	26.69	278	6	0.727	ng	# 1
39) Benzo(g,h,i)perylene	27.50	276	11	1.331	ng	# 1

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

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