

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091119\
 Data File : BN007588.D
 Acq On : 11 Sep 2019 15:17
 Operator : HP/JU
 Sample : K4698-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GW507-K2-35.0-40.0-090419

Quant Time: Sep 11 15:50:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 11 14:14:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	161	0.40	ng	0.00
7) Naphthalene-d8	10.49	136	601	0.40	ng	0.00
13) Acenaphthene-d10	14.34	164	351	0.40	ng	0.00
19) Phenanthrene-d10	17.09	188	612	0.40	ng	0.00
27) Chrysene-d12	21.28	240	1397	0.40	ng	0.00
34) Perylene-d12	23.51	264	703	0.40	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.35	112	86	0.18	ng	0.00
5) Phenol-d6	6.90	99	81	0.12	ng	0.00
8) Nitrobenzene-d5	8.87	82	193	0.53	ng	0.01
11) 2-Methylnaphthalene-d10	12.09	152	356	0.34	ng	0.00
14) 2,4,6-Tribromophenol	15.84	330	47	0.35	ng	0.00
15) 2-Fluorobiphenyl	12.97	172	512	0.35	ng	0.00
25) Fluoranthene-d10	19.12	212	708	0.36	ng	0.00
29) Terphenyl-d14	19.73	244	822	0.23	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) n-Nitrosodimethylamine	3.04	42	3174	33.002	ng	# 100
6) bis(2-Chloroethyl)ether	7.21	93	17	0.248	ng	# 64
9) Naphthalene	10.54	128	169	0.100	ng	# 1
10) Hexachlorobutadiene	10.85	225	7	0.021	ng	# 71
12) 2-Methylnaphthalene	12.16	142	104	0.088	ng	# 25
16) Acenaphthylene	14.06	152	62	0.033	ng	# 1
17) Acenaphthene	14.41	154	72	0.060	ng	# 71
18) Fluorene	15.39	166	136	0.087	ng	# 94
20) 4-Bromophenyl-phenylether	16.29	248	11	0.030	ng	# 1
21) Hexachlorobenzene	16.35	284	12	0.032	ng	# 41
22) Pentachlorophenol	16.74	266	46	0.535	ng	# 68
23) Phenanthrene	17.13	178	1051	0.556	ng	# 91
24) Anthracene	17.22	178	160	0.092	ng	# 73
26) Fluoranthene	19.15	202	962	0.420	ng	# 80
28) Pyrene	19.51	202	973	0.194	ng	# 94
30) Benzo(a)anthracene	21.26	228	394	0.079	ng	# 1
31) Chrysene	21.31	228	518	0.102	ng	# 1
32) Bis(2-ethylhexyl)phthalate	21.23	149	5599	3.273	ng	# 92
35) Benzo(b)fluoranthene	22.84	252	146	0.058	ng	# 1
36) Benzo(k)fluoranthene	22.90	252	738	0.299	ng	# 1
37) Benzo(a)pyrene	23.41	252	86	0.039	ng	# 1
38) Dibenzo(a,h)anthracene	25.75	278	62	0.028	ng	# 1
39) Benzo(g,h,i)perylene	26.41	276	158	0.070	ng	# 1

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

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