

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022120\
 Data File : BN009806.D
 Acq On : 21 Feb 2020 13:57
 Operator : CG/JU
 Sample : PB126934BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB126934BS

Quant Time: Feb 21 16:37:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN021920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 19 15:57:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	1117	0.40	ng	-0.02
7) Naphthalene-d8	10.16	136	3832	0.40	ng	-0.01
13) Acenaphthene-d10	14.04	164	2498	0.40	ng	-0.02
19) Phenanthrene-d10	16.80	188	5586	0.40	ng	-0.01
27) Chrysene-d12	21.02	240	4991	0.40	ng	-0.02
34) Perylene-d12	23.12	264	5433	0.40	ng	-0.03

System Monitoring Compounds						
4) 2-Fluorophenol	5.07	112	1002	0.40	ng	-0.02
5) Phenol-d6	6.64	99	1072	0.35	ng	0.00
8) Nitrobenzene-d5	8.58	82	1080	0.34	ng	-0.01
11) 2-Methylnaphthalene-d10	11.77	152	2897	0.39	ng	-0.02
14) 2,4,6-Tribromophenol	15.57	330	362	0.38	ng	-0.02
15) 2-Fluorobiphenyl	12.67	172	3615	0.38	ng	-0.02
25) Fluoranthene-d10	18.84	212	7134	0.37	ng	-0.02
29) Terphenyl-d14	19.46	244	4501	0.38	ng	-0.01

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.08	88	471	0.412	ng	95
3) n-Nitrosodimethylamine	3.43	42	670	0.357	ng	# 90
6) bis(2-Chloroethyl)ether	6.87	93	1182	0.402	ng	89
9) Naphthalene	10.20	128	4061	0.389	ng	100
10) Hexachlorobutadiene	10.48	225	910	0.396	ng	# 95
12) 2-Methylnaphthalene	11.85	142	2796	0.390	ng	98
16) Acenaphthylene	13.75	152	4059	0.381	ng	100
17) Acenaphthene	14.11	154	2895	0.387	ng	99
18) Fluorene	15.10	166	3723	0.377	ng	98
20) 4-Bromophenyl-phenylether	16.01	248	693	0.097	ng	# 85
21) Hexachlorobenzene	16.12	284	1390	0.400	ng	98
22) Pentachlorophenol	16.50	266	267	0.408	ng	92
23) Phenanthrene	16.84	178	6175	0.371	ng	99
24) Anthracene	16.94	178	5138	0.366	ng	99
26) Fluoranthene	18.88	202	7293	0.366	ng	99
28) Pyrene	19.24	202	7478	0.395	ng	99
30) Benzo(a)anthracene	21.01	228	5891	0.356	ng	100
31) Chrysene	21.06	228	7459	0.389	ng	98
32) Bis(2-ethylhexyl)phthalate	20.94	149	3439	0.426	ng	100
33) Indeno(1,2,3-cd)pyrene	25.17	276	7718	0.387	ng	98
35) Benzo(b)fluoranthene	22.50	252	6582	0.331	ng	99
36) Benzo(k)fluoranthene	22.55	252	8039	0.372	ng	99
37) Benzo(a)pyrene	23.03	252	6811	0.376	ng	# 87
38) Dibenzo(a,h)anthracene	25.18	278	5984	0.335	ng	99
39) Benzo(g,h,i)perylene	25.79	276	6633	0.349	ng	99

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

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