

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121118\
 Data File : BE098337.D
 Acq On : 11 Dec 2018 17:43
 Operator : SJ/JU
 Sample : J6259-08MSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 BPS1-TT-MW306I-20181205MSD

Manual Integrations
 APPROVED

Sohil
 12/12/2018 5:28:36 PM

Quant Time: Dec 12 11:40:34 2018

Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE121018.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Mon Dec 10 14:59:55 2018

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	1383	0.40	ng	0.00
7) Naphthalene-d8	10.67	136	5592	0.40	ng	0.00
13) Acenaphthene-d10	14.53	164	3375	0.40	ng	0.00
19) Phenanthrene-d10	17.27	188	8582	0.40	ng	-0.01
27) Chrysene-d12	21.46	240	10588	0.40	ng	0.00
34) Perylene-d12	24.01	264	10112	0.40	ng	0.00

System Monitoring Compounds

4) 2-Fluorophenol	5.46	112	600	0.19	ng	0.01
5) Phenol-d6	7.05	99	501	0.11	ng	0.00
8) Nitrobenzene-d5	9.03	82	2196	0.43	ng	0.01
11) 2-Methylnaphthalene-d10	12.27	152	3455m	0.38	ng	0.00
14) 2,4,6-Tribromophenol	16.02	330	559	0.45	ng	0.00
15) 2-Fluorobiphenyl	13.15	172	7391	0.52	ng	0.00
25) Fluoranthene-d10	19.31	212	49768	0.45	ng	0.00
29) Terphenyl-d14	19.90	244	12250	0.48	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.37	88	2165	0.905	ng	# 90
3) n-Nitrosodimethylamine	3.69	42	244	0.113	ng	# 100
6) bis(2-Chloroethyl)ether	7.29	93	1517	0.348	ng	88
9) Naphthalene	10.71	128	24450	0.434	ng	100
10) Hexachlorobutadiene	10.99	225	1323	0.369	ng	98
12) 2-Methylnaphthalene	12.34	142	4197	0.459	ng	96
16) Acenaphthylene	14.24	152	6788	0.429	ng	100
17) Acenaphthene	14.59	154	3869	0.407	ng	99
18) Fluorene	15.57	166	5441	0.428	ng	98
20) 4-Bromophenyl-phenylether	16.47	248	1867	0.404	ng	# 81
21) Hexachlorobenzene	16.58	284	1954	0.393	ng	94
22) Pentachlorophenol	16.93	266	689	0.765	ng	94
23) Phenanthrene	17.32	178	9269	0.444	ng	99
24) Anthracene	17.41	178	8269	0.445	ng	99
26) Fluoranthene	19.34	202	12816	0.464	ng	100
28) Pyrene	19.70	202	13184	0.397	ng	99
30) Benzo(a)anthracene	21.44	228	13868	0.460	ng	98
31) Chrysene	21.50	228	13706	0.434	ng	99
32) Bis(2-ethylhexyl)phthalate	21.37	149	42554	0.695	ng	99
33) Indeno(1,2,3-cd)pyrene	26.70	276	14027	0.446	ng	98
35) Benzo(b)fluoranthene	23.22	252	12451	0.450	ng	99
36) Benzo(k)fluoranthene	23.28	252	14711	0.457	ng	98
37) Benzo(a)pyrene	23.89	252	12825	0.449	ng	98
38) Dibenzo(a,h)anthracene	26.72	278	11750	0.437	ng	96
39) Benzo(g,h,i)perylene	27.52	276	12035	0.444	ng	98

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

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