

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN032319\
 Data File : BN004920.D
 Acq On : 23 Mar 2019 14:58
 Operator : JU/SJ
 Sample : PB118204BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB118204BS

Manual Integrations
 APPROVED

Sohil
 3/25/2019 1:34:20 PM

Quant Time: Mar 24 01:33:15 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN032019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 20 16:06:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	1109	0.40	ng	0.00
7) Naphthalene-d8	10.48	136	4250	0.40	ng	-0.01
13) Acenaphthene-d10	14.33	164	2441	0.40	ng	-0.01
19) Phenanthrene-d10	17.08	188	5822	0.40	ng	-0.01
27) Chrysene-d12	21.28	240	8083	0.40	ng	-0.01
34) Perylene-d12	23.53	264	9888	0.40	ng	-0.01

System Monitoring Compounds						
4) 2-Fluorophenol	5.28	112	1134	0.38	ng	0.00
5) Phenol-d6	6.87	99	1322	0.35	ng	0.00
8) Nitrobenzene-d5	8.86	82	1023	0.36	ng	-0.01
11) 2-Methylnaphthalene-d10	12.07	152	2125	0.29	ng	0.00
14) 2,4,6-Tribromophenol	15.84	330	533	0.32	ng	-0.01
15) 2-Fluorobiphenyl	12.95	172	3204	0.32	ng	-0.01
25) Fluoranthene-d10	19.12	212	6240	0.32	ng	0.00
29) Terphenyl-d14	19.72	244	5142	0.31	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.22	88	328	0.260	ng	# 69
3) n-Nitrosodimethylamine	3.55	42	271	0.364	ng	# 87
6) bis(2-Chloroethyl)ether	7.12	93	1082	0.320	ng	98
9) Naphthalene	10.53	128	3792	0.318	ng	96
10) Hexachlorobutadiene	10.80	225	674	0.296	ng	# 98
12) 2-Methylnaphthalene	12.15	142	2677	0.302	ng	98
16) Acenaphthylene	14.05	152	3832	0.305	ng	99
17) Acenaphthene	14.39	154	2461	0.299	ng	98
18) Fluorene	15.38	166	3323	0.289	ng	99
20) 4-Bromophenyl-phenylether	16.27	248	1155	0.304	ng	96
21) Hexachlorobenzene	16.39	284	1206	0.277	ng	94
22) Pentachlorophenol	16.75	266	714m	0.839	ng	
23) Phenanthrene	17.12	178	5794	0.308	ng	99
24) Anthracene	17.22	178	5284	0.315	ng	99
26) Fluoranthene	19.15	202	7441	0.308	ng	99
28) Pyrene	19.52	202	7737	0.314	ng	99
30) Benzo(a)anthracene	21.27	228	8749	0.330	ng	98
31) Chrysene	21.32	228	8640	0.309	ng	98
32) Bis(2-ethylhexyl)phthalate	21.17	149	4640	0.438	ng	97
33) Indeno(1,2,3-cd)pyrene	25.80	276	14768	0.366	ng	98
35) Benzo(b)fluoranthene	22.85	252	10378	0.290	ng	# 90
36) Benzo(k)fluoranthene	22.90	252	10284	0.292	ng	# 90
37) Benzo(a)pyrene	23.43	252	10307	0.314	ng	# 89
38) Dibenzo(a,h)anthracene	25.81	278	12392	0.326	ng	# 94
39) Benzo(g,h,i)perylene	26.50	276	12920	0.334	ng	96

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

