

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082119\
 Data File : BN007299.D
 Acq On : 21 Aug 2019 20:34
 Operator : HP/JU
 Sample : K4440-11
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 S303-J-1.5-2.0-082019

Manual Integrations
 APPROVED

Jagrut
 8/26/2019 12:24:46 PM

Quant Time: Aug 22 05:42:54 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082019.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Tue Aug 20 18:52:43 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.41	152	6025	0.40	ng	0.00
7) Naphthalene-d8	10.16	136	22626	0.40	ng	0.00
13) Acenaphthene-d10	14.06	164	12418	0.40	ng	0.00
19) Phenanthrene-d10	16.80	188	28450	0.40	ng	0.00
27) Chrysene-d12	21.03	240	29790	0.40	ng	0.00
34) Perylene-d12	23.14	264	34770	0.40	ng	0.00

System Monitoring Compounds

4) 2-Fluorophenol	5.04	112	3190	0.17	ng	0.00
5) Phenol-d6	6.61	99	4019	0.18	ng	0.00
8) Nitrobenzene-d5	8.54	82	2817	0.16	ng	0.00
11) 2-Methylnaphthalene-d10	11.77	152	6145	0.16	ng	0.00
14) 2,4,6-Tribromophenol	15.56	330	1218	0.17	ng	0.00
15) 2-Fluorobiphenyl	12.68	172	7898	0.15	ng	0.00
25) Fluoranthene-d10	18.85	212	14484	0.16	ng	0.00
29) Terphenyl-d14	19.48	244	10849	0.15	ng	0.00

Target Compounds

						Qvalue
16) Acenaphthylene	13.77	152	3057	0.046	ng	98
17) Acenaphthene	14.12	154	2734	0.065	ng	99
18) Fluorene	15.12	166	3653	0.067	ng	94
23) Phenanthrene	16.85	178	58777	0.659	ng	100
24) Anthracene	16.94	178	16096	0.200	ng	# 94
26) Fluoranthene	18.88	202	289022	2.676	ng	99
28) Pyrene	19.25	202	258564	2.321	ng	# 94
30) Benzo(a)anthracene	21.02	228	192989	1.742	ng	95
31) Chrysene	21.07	228	158256m	1.400	ng	
32) Bis(2-ethylhexyl)phthalate	20.98	149	8021	0.181	ng	# 76
33) Indeno(1,2,3-cd)pyrene	25.21	276	114083	0.839	ng	95
35) Benzo(b)fluoranthene	22.52	252	234401	1.808	ng	# 94
36) Benzo(k)fluoranthene	22.56	252	90824m	0.715	ng	
37) Benzo(a)pyrene	23.05	252	181312	1.552	ng	# 94
38) Dibenzo(a,h)anthracene	25.22	278	30629	0.250	ng	# 58
39) Benzo(g,h,i)perylene	25.84	276	112105	0.904	ng	93

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

