

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092519\
 Data File : BN007864.D
 Acq On : 25 Sep 2019 21:16
 Operator : JU
 Sample : K4994-03 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EXT-SW092-091819-1

Manual Integrations
 APPROVED

mohammad
 9/27/2019 8:28:06 AM

Quant Time: Sep 26 09:07:58 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091819.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Sep 18 18:24:40 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	10416	0.40	ng	0.00
7) Naphthalene-d8	10.46	136	38791	0.40	ng	-0.01
13) Acenaphthene-d10	14.31	164	22117	0.40	ng	-0.01
19) Phenanthrene-d10	17.06	188	41660	0.40	ng	-0.01
27) Chrysene-d12	21.25	240	46135	0.40	ng	-0.02
34) Perylene-d12	23.47	264	56966	0.40	ng	-0.02

System Monitoring Compounds

4) 2-Fluorophenol	5.31	112	1315	0.04	ng	0.00
5) Phenol-d6	6.87	99	1705	0.03	ng	0.00
8) Nitrobenzene-d5	8.83	82	1443	0.04	ng	0.00
11) 2-Methylnaphthalene-d10	12.05	152	2730	0.04	ng	-0.01
14) 2,4,6-Tribromophenol	15.81	330	397	0.03	ng	-0.01
15) 2-Fluorobiphenyl	12.93	172	3909	0.04	ng	-0.01
25) Fluoranthene-d10	19.09	212	5439m	0.04	ng	-0.01
29) Terphenyl-d14	19.69	244	7518	0.07	ng	-0.02

Target Compounds

						Qvalue
9) Naphthalene	10.51	128	38510	0.354	ng	100
12) 2-Methylnaphthalene	12.13	142	32354	0.440	ng	100
16) Acenaphthylene	14.03	152	250716	2.137	ng	99
17) Acenaphthene	14.38	154	34111	0.451	ng	# 88
18) Fluorene	15.36	166	74491	0.769	ng	86
23) Phenanthrene	17.10	178	1786091	13.763	ng	100
24) Anthracene	17.19	178	349401	2.981	ng	98
26) Fluoranthene	19.12	202	2583619	16.010	ng	98
28) Pyrene	19.49	202	3789318	24.083	ng	# 95
30) Benzo(a)anthracene	21.24	228	1970175	11.597	ng	96
31) Chrysene	21.29	228	2094542	12.650	ng	97
32) Bis(2-ethylhexyl)phthalate	21.19	149	26621m	0.305	ng	
33) Indeno(1,2,3-cd)pyrene	25.68	276	1004473	4.845	ng	97
35) Benzo(b)fluoranthene	22.81	252	2015325m	10.190	ng	
36) Benzo(k)fluoranthene	22.85	252	575663m	2.944	ng	
37) Benzo(a)pyrene	23.37	252	1680567	9.043	ng	98
38) Dibenzo(a,h)anthracene	25.68	278	312816	1.645	ng	93
39) Benzo(a,h,i)perylene	26.35	276	1199055	6.372	ng	99

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

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