

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE062719\
 Data File : BE100217.D
 Acq On : 27 Jun 2019 22:42
 Operator : JU/SJ
 Sample : K3428-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 EXT-013-SW159-061319

Manual Integrations
 APPROVED

mohammad
 7/1/2019 8:12:21 AM

Quant Time: Jun 28 05:57:59 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE062519.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Jun 28 05:41:07 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	1028	0.40	ng	-0.01
7) Naphthalene-d8	10.45	136	3583	0.40	ng	0.00
13) Acenaphthene-d10	14.33	164	2778	0.40	ng	0.00
19) Phenanthrene-d10	17.07	188	7754	0.40	ng	0.00
27) Chrysene-d12	21.24	240	9654	0.40	ng	0.00
34) Perylene-d12	23.67	264	11367	0.40	ng	0.00

System Monitoring Compounds

4) 2-Fluorophenol	5.23	112	869	0.38	ng	0.00
5) Phenol-d6	6.83	99	1218	0.39	ng	-0.02
8) Nitrobenzene-d5	8.82	82	1224	0.34	ng	-0.01
11) 2-Methylnaphthalene-d10	12.05	152	2541	0.38	ng	-0.02
14) 2,4,6-Tribromophenol	15.81	330	687	0.43	ng	-0.03
15) 2-Fluorobiphenyl	12.94	172	3632	0.33	ng	-0.02
25) Fluoranthene-d10	19.10	212	38472	0.34	ng	0.00
29) Terphenyl-d14	19.70	244	7210	0.38	ng	0.00

Target Compounds

						Qvalue
9) Naphthalene	10.50	128	2787	0.071	ng	# 95
12) 2-Methylnaphthalene	12.12	142	560	0.089	ng	93
16) Acenaphthylene	14.04	152	3310	0.244	ng	97
17) Acenaphthene	14.38	154	737	0.097	ng	97
18) Fluorene	15.37	166	1654	0.151	ng	# 83
23) Phenanthrene	17.11	178	41080	2.237	ng	99
24) Anthracene	17.20	178	4178	0.269	ng	97
26) Fluoranthene	19.13	202	64706	2.166	ng	99
28) Pyrene	19.49	202	81420	3.293	ng	# 96
30) Benzo(a)anthracene	21.23	228	46641	1.639	ng	94
31) Chrysene	21.28	228	45932	1.376	ng	95
32) Bis(2-ethylhexyl)phthalate	21.16	149	16968	0.519	ng	100
33) Indeno(1,2,3-cd)pyrene	26.23	276	21398	0.450	ng	# 95
35) Benzo(b)fluoranthene	22.93	252	45158	1.477	ng	93
36) Benzo(k)fluoranthene	22.97	252	12180m	0.333	ng	
37) Benzo(a)pyrene	23.56	252	33024m	1.042	ng	
38) Dibenzo(a,h)anthracene	26.24	278	6390	0.189	ng	# 75
39) Benzo(a,h,i)perylene	27.02	276	23636	0.668	ng	96

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE062719\
 Data File : BE100217.D
 Acq On : 27 Jun 2019 22:42
 Operator : JU/SJ
 Sample : K3428-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 EXT-013-SW159-061319

Manual Integrations
 APPROVED

mohammad
 7/1/2019 8:12:21 AM

Quant Time: Jun 28 05:57:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE062519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 28 05:41:07 2019
 Response via : Initial Calibration

