

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE070219\
 Data File : BE100315.D
 Acq On : 2 Jul 2019 22:53
 Operator : JU/SJ
 Sample : K3504-17 10X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PST-013-SW105-062019

Manual Integrations
 APPROVED

mohammad
 7/3/2019 10:49:19 PM

Quant Time: Jul 03 04:01:03 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE062819.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Jun 28 14:06:53 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	806	0.40	ng	0.00
7) Naphthalene-d8	10.43	136	3019	0.40	ng	-0.01
13) Acenaphthene-d10	14.33	164	2243	0.40	ng	0.01
19) Phenanthrene-d10	17.07	188	6333	0.40	ng	0.01
27) Chrysene-d12	21.25	240	8332	0.40	ng	0.00
34) Perylene-d12	23.68	264	9786	0.40	ng	0.01

System Monitoring Compounds

4) 2-Fluorophenol	5.23	112	58	0.03	ng	0.01
5) Phenol-d6	6.85	99	64	0.02	ng	0.01
8) Nitrobenzene-d5	8.82	82	68	0.02	ng	0.00
11) 2-Methylnaphthalene-d10	12.05	152	139	0.02	ng	0.00
14) 2,4,6-Tribromophenol	15.82	330	65	0.04	ng	0.00
15) 2-Fluorobiphenyl	12.94	172	197	0.02	ng	0.00
25) Fluoranthene-d10	19.10	212	1437	0.02	ng	0.00
29) Terphenyl-d14	19.70	244	814	0.05	ng	0.00

Target Compounds

						Qvalue
9) Naphthalene	10.48	128	22784	0.684	ng	98
12) 2-Methylnaphthalene	12.12	142	3452	0.591	ng	97
16) Acenaphthylene	14.04	152	3226	0.300	ng	# 94
17) Acenaphthene	14.38	154	8425	1.381	ng	98
18) Fluorene	15.35	166	11894	1.319	ng	95
23) Phenanthrene	17.11	178	256266	16.696	ng	100
24) Anthracene	17.19	178	56786	4.108	ng	# 93
26) Fluoranthene	19.13	202	434494	17.772	ng	99
28) Pyrene	19.49	202	412700	18.709	ng	# 95
30) Benzo(a)anthracene	21.23	228	262111	9.568	ng	98
31) Chrysene	21.28	228	207765	7.552	ng	97
33) Indeno(1,2,3-cd)pyrene	26.24	276	133761	3.676	ng	# 97
35) Benzo(b)fluoranthene	22.93	252	251532	9.480	ng	# 96
36) Benzo(k)fluoranthene	22.98	252	107607m	3.633	ng	
37) Benzo(a)pyrene	23.57	252	204642	7.696	ng	# 93
38) Dibenzo(a,h)anthracene	26.25	278	33250	1.197	ng	93
39) Benzo(a,h,i)perylene	27.04	276	135817	4.760	ng	99

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE070219\
Data File : BE100315.D
Acq On : 2 Jul 2019 22:53
Operator : JU/SJ
Sample : K3504-17 10X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
PST-013-SW105-062019

Manual Integrations
APPROVED

mohammad
7/3/2019 10:49:19 PM

Quant Time: Jul 03 04:01:03 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE062819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 28 14:06:53 2019
Response via : Initial Calibration

