

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE070119\
 Data File : BE100283.D
 Acq On : 2 Jul 2019 2:56
 Operator : JU/SJ
 Sample : K3446-15
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 EXT-008-SW114-061819

Manual Integrations
 APPROVED

mohammad
 7/3/2019 10:52:50 PM

Quant Time: Jul 02 04:14:50 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	7.64	152	1099	0.40	ng	0.00
7) Naphthalene-d8	10.43	136	3758	0.40	ng	-0.01
13) Acenaphthene-d10	14.31	164	2820	0.40	ng	0.00
19) Phenanthrene-d10	17.05	188	7556	0.40	ng	0.00
27) Chrysene-d12	21.24	240	10402	0.40	ng	0.00
34) Perylene-d12	23.68	264	12001	0.40	ng	0.02

System Monitoring Compounds

4) 2-Fluorophenol	5.23	112	809	0.29	ng	0.01
5) Phenol-d6	6.83	99	1137	0.31	ng	-0.01
8) Nitrobenzene-d5	8.81	82	1165	0.31	ng	-0.01
11) 2-Methylnaphthalene-d10	12.04	152	2536	0.36	ng	-0.01
14) 2,4,6-Tribromophenol	15.81	330	635	0.32	ng	-0.01
15) 2-Fluorobiphenyl	12.93	172	3432	0.31	ng	-0.01
25) Fluoranthene-d10	19.10	212	35917	0.33	ng	0.00
29) Terphenyl-d14	19.70	244	7806	0.36	ng	0.00

Target Compounds

						Qvalue
9) Naphthalene	10.48	128	13257	0.320	ng	99
12) 2-Methylnaphthalene	12.11	142	1945	0.267	ng	98
16) Acenaphthylene	14.04	152	13908	1.028	ng	99
17) Acenaphthene	14.38	154	1483	0.193	ng	97
18) Fluorene	15.35	166	3616	0.319	ng	# 92
23) Phenanthrene	17.09	178	105500	5.761	ng	100
24) Anthracene	17.19	178	18265	1.108	ng	# 96
26) Fluoranthene	19.13	202	334154	11.455	ng	99
28) Pyrene	19.50	202	339926	12.343	ng	# 95
30) Benzo(a)anthracene	21.23	228	246697	7.213	ng	98
31) Chrysene	21.28	228	241210	7.023	ng	95
32) Bis(2-ethylhexyl)phthalate	21.16	149	115726	2.336	ng	100
33) Indeno(1,2,3-cd)pyrene	26.25	276	153905	3.388	ng	# 96
35) Benzo(b)fluoranthene	22.94	252	310144	9.532	ng	# 97
36) Benzo(k)fluoranthene	22.98	252	100368m	2.763	ng	
37) Benzo(a)pyrene	23.57	252	215940m	6.622	ng	
38) Dibenzo(a,h)anthracene	26.26	278	40548	1.191	ng	# 91
39) Benzo(a,h,i)perylene	27.04	276	156357	4.469	ng	98

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

