

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE043019\
 Data File : BE099439.D
 Acq On : 30 Apr 2019 23:25
 Operator : JU/SJ
 Sample : K2568-16
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PST-003-B175-042319

Manual Integrations
 APPROVED

Sohil
 5/1/2019 12:02:50 PM

Quant Time: May 01 02:28:09 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE042919.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Mon Apr 29 18:36:43 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	2809	0.40	ng	0.00
7) Naphthalene-d8	10.60	136	10191	0.40	ng	0.00
13) Acenaphthene-d10	14.46	164	7532	0.40	ng	0.00
19) Phenanthrene-d10	17.19	188	24519	0.40	ng	0.00
27) Chrysene-d12	21.37	240	33806	0.40	ng	0.00
34) Perylene-d12	23.86	264	38600	0.40	ng	0.00

System Monitoring Compounds

4) 2-Fluorophenol	5.39	112	3011	0.39	ng	0.00
5) Phenol-d6	6.97	99	4413	0.43	ng	0.00
8) Nitrobenzene-d5	8.96	82	3675	0.38	ng	0.00
11) 2-Methylnaphthalene-d10	12.19	152	6607	0.36	ng	0.00
14) 2,4,6-Tribromophenol	15.93	330	2071	0.44	ng	-0.01
15) 2-Fluorobiphenyl	13.07	172	9671	0.34	ng	-0.02
25) Fluoranthene-d10	19.22	212	111823	0.34	ng	0.00
29) Terphenyl-d14	19.81	244	23794	0.38	ng	0.00

Target Compounds

						Qvalue
9) Naphthalene	10.65	128	12873	0.116	ng	# 99
12) 2-Methylnaphthalene	12.27	142	786	0.041	ng	# 94
16) Acenaphthylene	14.18	152	3624	0.098	ng	99
17) Acenaphthene	14.51	154	845	0.040	ng	96
18) Fluorene	15.49	166	1567	0.050	ng	# 93
23) Phenanthrene	17.23	178	57785	0.913	ng	99
24) Anthracene	17.32	178	6205	0.103	ng	# 84
26) Fluoranthene	19.25	202	161493	1.704	ng	100
28) Pyrene	19.61	202	156644	1.778	ng	# 96
30) Benzo(a)anthracene	21.34	228	97177	0.902	ng	97
31) Chrysene	21.40	228	107492	1.020	ng	97
32) Bis(2-ethylhexyl)phthalate	21.26	149	17159	0.041	ng	99
33) Indeno(1,2,3-cd)pyrene	26.50	276	63420	0.509	ng	97
35) Benzo(b)fluoranthene	23.09	252	128948	1.159	ng	# 96
36) Benzo(k)fluoranthene	23.14	252	48306m	0.450	ng	
37) Benzo(a)pyrene	23.74	252	88804	0.844	ng	99
38) Dibenzo(a,h)anthracene	26.52	278	16491	0.153	ng	# 81
39) Benzo(g,h,i)perylene	27.32	276	62625	0.581	ng	99

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

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