

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE050219\
 Data File : BE099522.D
 Acq On : 3 May 2019 9:07
 Operator : JU/SJ
 Sample : K2589-14DL 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PST-001-B188-042419DL

Manual Integrations
 APPROVED

Sohil
 5/3/2019 4:24:26 PM

Quant Time: May 03 13:38:37 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	7.80	152	2808	0.40	ng	-0.01
7) Naphthalene-d8	10.60	136	10937	0.40	ng	0.00
13) Acenaphthene-d10	14.46	164	9200	0.40	ng	0.00
19) Phenanthrene-d10	17.19	188	31302	0.40	ng	0.00
27) Chrysene-d12	21.37	240	38063	0.40	ng	0.00
34) Perylene-d12	23.87	264	44397	0.40	ng	0.00

System Monitoring Compounds

4) 2-Fluorophenol	5.39	112	346	0.04	ng	0.00
5) Phenol-d6	6.97	99	476	0.05	ng	0.00
8) Nitrobenzene-d5	8.96	82	392	0.04	ng	0.00
11) 2-Methylnaphthalene-d10	12.19	152	598	0.03	ng	0.00
14) 2,4,6-Tribromophenol	15.94	330	262	0.05	ng	0.00
15) 2-Fluorobiphenyl	13.09	172	1150	0.03	ng	0.00
25) Fluoranthene-d10	19.22	212	10766	0.03	ng	0.00
29) Terphenyl-d14	19.82	244	3323	0.05	ng	0.00

Target Compounds

						Qvalue
9) Naphthalene	10.65	128	11283	0.095	ng	# 98
12) 2-Methylnaphthalene	12.27	142	1452	0.070	ng	95
16) Acenaphthylene	14.18	152	1362	0.030	ng	# 91
17) Acenaphthene	14.51	154	4286	0.167	ng	98
18) Fluorene	15.49	166	5912	0.156	ng	# 96
22) Pentachlorophenol	16.84	266	159	0.231	ng	# 76
23) Phenanthrene	17.23	178	131345	1.626	ng	100
24) Anthracene	17.32	178	24810	0.322	ng	# 94
26) Fluoranthene	19.25	202	166444	1.376	ng	99
28) Pyrene	19.62	202	146941	1.481	ng	# 95
30) Benzo(a)anthracene	21.35	228	87477	0.721	ng	99
31) Chrysene	21.40	228	70680	0.596	ng	96
33) Indeno(1,2,3-cd)pyrene	26.51	276	32814	0.234	ng	97
35) Benzo(b)fluoranthene	23.10	252	74293	0.580	ng	97
36) Benzo(k)fluoranthene	23.15	252	27845m	0.226	ng	
37) Benzo(a)pyrene	23.76	252	55694	0.460	ng	99
38) Dibenzo(a,h)anthracene	26.54	278	9300	0.075	ng	# 83
39) Benzo(g,h,i)perylene	27.34	276	30544	0.246	ng	97

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

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