Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE041519\
Data File : BE099327.D
Acq On : 15 Apr 2019 15:53
Operator : JU/SJ
Sample : K2366-03MSD
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
SA-PZ118I-20190409MSD

Manual Integrations APPROVED
Sohil
4/23/2019 3:45:43 AM

Quant Time: Apr 16 03:55:40 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE040219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 02 17:16:42 2019
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	3200	0.40	ng		-0.02
7) Naphthalene-d8	10.61	136	11407	0.40	ng		-0.01
13) Acenaphthene-d10	14.46	164	7119	0.40	ng		-0.03
19) Phenanthrene-d10	17.19	188	20382	0.40	ng		-0.03
27) Chrysene-d12	21.37	240	32205	0.40	ng		-0.02
34) Perylene-d12	23.87	264	32519	0.40	ng		-0.03
System Monitoring Compounds							
4) 2-Fluorophenol	5.39	112	1587	0.19	ng		-0.02
5) Phenol-d6	6.98	99	1191	0.10	ng		-0.01
8) Nitrobenzene-d5	8.96	82	3756	0.48	ng		-0.03
11) 2-Methylnaphthalene-d10	12.21	152	7833m	0.39	ng		-0.01
14) 2,4,6-Tribromophenol	15.94	330	1224	0.37	ng		-0.01
15) 2-Fluorobiphenyl	13.09	172	13683	0.49	ng		-0.01
25) Fluoranthene-d10	19.22	212	120744	0.45	ng		-0.02
29) Terphenyl-d14	19.82	244	26204	0.45	ng		-0.02
Target Compounds							Qvalue
2) 1,4-Dioxane	3.33	88	1589	0.364	ng	#	71
3) n-Nitrosodimethylamine	3.65	42	272	0.106	ng	#	33
6) bis(2-Chloroethyl)ether	7.24	93	4376	0.417	ng		97
9) Naphthalene	10.65	128	53158	0.428	ng		100
10) Hexachlorobutadiene	10.94	225	2900	0.360	ng		97
12) 2-Methylnaphthalene	12.28	142	8787	0.416	ng		98
16) Acenaphthylene	14.18	152	12953	0.384	ng		99
17) Acenaphthene	14.53	154	8142	0.409	ng		100
18) Fluorene	15.50	166	11679	0.406	ng		98
20) 4-Bromophenyl-phenylether	16.38	248	4655	0.402	ng	#	77
21) Hexachlorobenzene	16.49	284	4414	0.401	ng		97
22) Pentachlorophenol	16.84	266	3330	0.728	ng		93
23) Phenanthrene	17.23	178	22882	0.434	ng		99
24) Anthracene	17.32	178	20661	0.415	ng		99
26) Fluoranthene	19.25	202	34495	0.443	ng		99
28) Pyrene	19.61	202	37453	0.440	ng		99
30) Benzo(a)anthracene	21.34	228	43179	0.426	ng		97
31) Chrysene	21.40	228	43121	0.434	ng		98
32) Bis(2-ethylhexyl)phthalate	21.27	149	57982	0.449	ng		100
33) Indeno(1,2,3-cd)pyrene	26.51	276	44819	0.394	ng		98
35) Benzo(b)fluoranthene	23.09	252	42727m	0.440	ng		
36) Benzo(k)fluoranthene	23.14	252	43720	0.463	ng		100
37) Benzo(a)pyrene	23.75	252	40288	0.442	ng	#	93
38) Dibenzo(a,h)anthracene	26.53	278	37541	0.414	ng		99
39) Benzo(g,h,i)perylene	27.32	276	37951	0.420	ng		100

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

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