

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070119\
 Data File : BN006671.D
 Acq On : 01 Jul 2019 22:49
 Operator : HP/JU
 Sample : K3504-20
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EXT-009-SW136-062019

Manual Integrations
 APPROVED

mohammad
 7/2/2019 2:52:40 PM

Quant Time: Jul 02 05:47:02 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN062819.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Jun 28 13:47:56 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	4645	0.40	ng	0.00
7) Naphthalene-d8	10.39	136	18942	0.40	ng	-0.01
13) Acenaphthene-d10	14.27	164	10533	0.40	ng	-0.01
19) Phenanthrene-d10	17.03	188	22183	0.40	ng	-0.01
27) Chrysene-d12	21.25	240	19900	0.40	ng	-0.02
34) Perylene-d12	23.48	264	24065	0.40	ng	-0.04

System Monitoring Compounds

4) 2-Fluorophenol	5.24	112	3333	0.29	ng	0.00
5) Phenol-d6	6.82	99	4658	0.31	ng	0.00
8) Nitrobenzene-d5	8.79	82	3245	0.25	ng	0.00
11) 2-Methylnaphthalene-d10	11.99	152	7721	0.26	ng	0.00
14) 2,4,6-Tribromophenol	15.78	330	1498	0.30	ng	-0.01
15) 2-Fluorobiphenyl	12.88	172	10330	0.26	ng	-0.01
25) Fluoranthene-d10	19.07	212	15516	0.24	ng	-0.01
29) Terphenyl-d14	19.67	244	11394	0.27	ng	-0.01

Target Compounds

						Qvalue
9) Naphthalene	10.44	128	19064	0.383	ng	99
12) 2-Methylnaphthalene	12.07	142	16151	0.470	ng	99
16) Acenaphthylene	13.99	152	163674	3.259	ng	100
17) Acenaphthene	14.33	154	11448	0.346	ng	95
18) Fluorene	15.32	166	36003	0.863	ng	# 87
22) Pentachlorophenol	16.72	266	330	0.285	ng	96
23) Phenanthrene	17.07	178	599846	9.510	ng	100
24) Anthracene	17.16	178	130250	2.380	ng	99
26) Fluoranthene	19.10	202	869715	11.346	ng	98
28) Pvrene	19.47	202	968668	13.911	ng	97
30) Benzo(a)anthracene	21.23	228	398306	6.122	ng	96
31) Chrysene	21.28	228	466457	6.651	ng	96
32) Bis(2-ethylhexyl)phthalate	21.15	149	130589	3.988	ng	99
33) Indeno(1,2,3-cd)pyrene	25.71	276	277061	3.903	ng	97
35) Benzo(b)fluoranthene	22.81	252	548411	6.699	ng	99
36) Benzo(k)fluoranthene	22.85	252	180402m	2.067	ng	
37) Benzo(a)pvrene	23.38	252	400165	5.229	ng	99
38) Dibenzo(a,h)anthracene	25.71	278	76802	1.149	ng	# 86
39) Benzo(g,h,i)perylene	26.40	276	294913	4.419	ng	100

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

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