

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082919\
 Data File : BN007517.D
 Acq On : 30 Aug 2019 00:05
 Operator : HP/JU
 Sample : K4514-02DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 S602-M-1.0-1.5-082319DL

Manual Integrations
 APPROVED

JAGRUT
 8/30/2019 8:58:06 AM

Quant Time: Aug 30 02:45:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 29 12:18:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	6179	0.40	ng	0.00
7) Naphthalene-d8	10.15	136	28057	0.40	ng	-0.01
13) Acenaphthene-d10	14.05	164	17341	0.40	ng	0.00
19) Phenanthrene-d10	16.79	188	40155	0.40	ng	0.00
27) Chrysene-d12	21.02	240	38891	0.40	ng	0.00
34) Perylene-d12	23.12	264	33184	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.03	112	1226	0.03	ng	0.00
5) Phenol-d6	6.59	99	1759	0.04	ng	0.00
8) Nitrobenzene-d5	8.53	82	1386	0.06	ng	0.00
11) 2-Methylnaphthalene-d10	11.76	152	2458	0.05	ng	0.00
14) 2,4,6-Tribromophenol	15.55	330	528	0.06	ng	0.00
15) 2-Fluorobiphenyl	12.66	172	3232	0.05	ng	0.00
25) Fluoranthene-d10	18.83	212	6352	0.05	ng	-0.01
29) Terphenyl-d14	19.46	244	4489	0.04	ng	-0.01
Target Compounds						
12) 2-Methylnaphthalene	11.83	142	2211	0.040	ng	Qvalue 97
17) Acenaphthene	14.10	154	6806	0.113	ng	99
18) Fluorene	15.10	166	5938	0.078	ng	98
23) Phenanthrene	16.84	178	96588	0.800	ng	99
24) Anthracene	16.92	178	18909	0.156	ng	97
26) Fluoranthene	18.86	202	320764	2.068	ng	99
28) Pyrene	19.23	202	280819	1.795	ng	# 95
30) Benzo(a)anthracene	20.99	228	157459	1.031	ng	99
31) Chrysene	21.05	228	153191	1.039	ng	99
32) Bis(2-ethylhexyl)phthalate	20.96	149	5457	0.036	ng	98
33) Indeno(1,2,3-cd)pyrene	25.17	276	79869	0.450	ng	99
35) Benzo(b)fluoranthene	22.50	252	191754	1.595	ng	# 96
36) Benzo(k)fluoranthene	22.53	252	68560m	0.577	ng	
37) Benzo(a)pyrene	23.02	252	125472m	1.097	ng	
38) Dibenzo(a,h)anthracene	25.18	278	21842	0.191	ng	# 90
39) Benzo(g,h,i)perylene	25.80	276	78124	0.687	ng	99

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

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