

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110719\
 Data File : BN008514.D
 Acq On : 08 Nov 2019 18:14
 Operator : JU
 Sample : K5699-05DL 2X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F2J02DL

Manual Integrations
 APPROVED

mohammad
 11/11/2019 8:09:16 AM

Quant Time: Nov 08 23:18:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 08 23:03:10 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	1657	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.31	136	7065	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.19	164	4257	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	9258	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	7246	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	6778	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	1412	0.13	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	4242	0.15	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	12204	0.607	ng/ul	98
5) 2-Methylnaphthalene	11.99	142	602	0.045	ng/ul	98
7) Acenaphthylene	13.90	152	2300	0.116	ng/ul	95
8) Acenaphthene	14.25	153	1490	0.093	ng/ul	99
9) Fluorene	15.24	166	1062	0.060	ng/ul#	97
12) Phenanthrene	16.97	178	2512	0.089	ng/ul	94
13) Anthracene	17.07	178	7256	0.294	ng/ul	100
15) Fluoranthene	19.00	202	16734	0.519	ng/ul	98
17) Pyrene	19.37	202	24258	0.686	ng/ul	99
18) Benzo(a)anthracene	21.13	228	10464	0.415	ng/ul	98
19) Chrysene	21.18	228	10619	0.331	ng/ul	99
21) Benzo(b)fluoranthene	22.67	252	25472m	1.008	ng/ul	
22) Benzo(k)fluoranthene	22.71	252	8372m	0.292	ng/ul	
23) Benzo(a)pyrene	23.21	252	9996	0.393	ng/ul#	74
24) Indeno(1,2,3-cd)pyrene	25.45	276	6288	0.218	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.45	278	1550	0.068	ng/ul#	78
26) Benzo(a,h,i)perylene	26.10	276	5257	0.203	ng/ul	97

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

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