

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111919\
 Data File : BN008648.D
 Acq On : 20 Nov 2019 02:28
 Operator : JU
 Sample : K5699-19
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F2J17

Manual Integrations
 APPROVED

mohammad
 11/21/2019 9:05:42 AM

Quant Time: Nov 20 04:53:19 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 20 04:45:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	3958	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	14704	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	8188	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	17044m	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	14347	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	13512m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	7104	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	17669	0.34	ng/ul	0.00
Target Compounds				Ovalue		
9) Fluorene	15.30	166	1147	0.029	ng/ul#	94
11) Pentachlorophenol	16.65	266	254	0.039	ng/ul	96
12) Phenanthrene	17.03	178	6252	0.109	ng/ul	99
13) Anthracene	17.12	178	4717	0.087	ng/ul	96
15) Fluoranthene	19.05	202	7985	0.113	ng/ul	91
17) Pyrene	19.41	202	7925m	0.138	ng/ul	
18) Benzo(a)anthracene	21.17	228	7772	0.135	ng/ul	98
19) Chrysene	21.22	228	7208	0.127	ng/ul	98
21) Benzo(b)fluoranthene	22.71	252	7893	0.142	ng/ul#	72
22) Benzo(k)fluoranthene	22.75	252	7646	0.138	ng/ul#	53
23) Benzo(a)pyrene	23.26	252	6123	0.117	ng/ul#	64
24) Indeno(1,2,3-cd)pyrene	25.49	276	8489	0.139	ng/ul	95
25) Dibenzo(a,h)anthracene	25.51	278	6614	0.134	ng/ul#	84
26) Benzo(g,h,i)perylene	26.15	276	7315	0.143	ng/ul#	85

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

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