

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111819\
 Data File : BN008609.D
 Acq On : 18 Nov 2019 18:19
 Operator : JU
 Sample : K5678-05DL 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLM69DL

Manual Integrations
 APPROVED

mohammad
 11/19/2019 2:45:01 PM

Quant Time: Nov 19 02:37:41 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 : Tue Nov 19 01:47:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	3804	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	13813	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	8354	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	17101	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	13159	0.40	ng/ul	0.00
20) Perylene-d12	23.38	264	14646	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	729	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	1740	0.03	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	3384	0.081	ng/ul#	92
5) 2-Methylnaphthalene	12.06	142	1281	0.044	ng/ul	100
7) Acenaphthylene	13.97	152	1151	0.025	ng/ul#	72
11) Pentachlorophenol	16.65	266	700	0.108	ng/ul	95
12) Phenanthrene	17.04	178	10335	0.179	ng/ul#	91
13) Anthracene	17.13	178	1805	0.033	ng/ul#	39
15) Fluoranthene	19.06	202	14018	0.198	ng/ul	86
17) Pyrene	19.42	202	27935	0.529	ng/ul	98
18) Benzo(a)anthracene	21.18	228	8492	0.160	ng/ul#	55
19) Chrysene	21.23	228	22712	0.437	ng/ul#	77
21) Benzo(b)fluoranthene	22.73	252	20665m	0.342	ng/ul	
22) Benzo(k)fluoranthene	22.77	252	4634m	0.077	ng/ul	
23) Benzo(a)pyrene	23.28	252	17684	0.312	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.53	276	10953	0.165	ng/ul#	63
25) Dibenzo(a,h)anthracene	25.54	278	3711	0.069	ng/ul#	1
26) Benzo(a,h,i)perylene	26.18	276	30398	0.548	ng/ul#	76

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

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