

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122819\
 Data File : BN009207.D
 Acq On : 27 Dec 2019 22:40
 Operator : JU
 Sample : K6276-15
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0C63

Manual Integrations
 APPROVED

mohammad
 12/29/2019 5:31:13 PM

Quant Time: Dec 27 23:33:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 27 23:08:48 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	2266	0.40	ng/ul	0.00
2) Naphthalene-d8	10.27	136	8113	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.14	164	4610	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.89	188	9680m	0.40	ng/ul	0.00
16) Chrysene-d12	21.09	240	8080	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	8068	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.86	152	3345	0.23	ng/ul	0.00
14) Fluoranthene-d10	18.93	212	8293	0.29	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	16.93	178	749	0.022	ng/ul#	77
15) Fluoranthene	18.95	202	2220	0.058	ng/ul	87
17) Pyrene	19.32	202	1864	0.045	ng/ul#	87
18) Benzo(a)anthracene	21.07	228	1212	0.034	ng/ul#	82
19) Chrysene	21.12	228	1550	0.043	ng/ul#	86
21) Benzo(b)fluoranthene	22.59	252	2775	0.074	ng/ul#	53
22) Benzo(k)fluoranthene	22.62	252	1282m	0.034	ng/ul	
23) Benzo(a)pyrene	23.12	252	1270	0.038	ng/ul#	20
24) Indeno(1,2,3-cd)pyrene	25.31	276	1166	0.032	ng/ul#	82
26) Benzo(g,h,i)perylene	25.94	276	1138	0.038	ng/ul#	66

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

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