

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122819\
 Data File : BN009241.D
 Acq On : 28 Dec 2019 22:52
 Operator : JU
 Sample : K6278-16DL 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0C69DL

Manual Integrations
 APPROVED

mohammad
 12/29/2019 5:33:16 PM

Quant Time: Dec 29 05:45:18 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 : Sun Dec 29 04:38:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	2627	0.40	ng/ul	0.00
2) Naphthalene-d8	10.26	136	9377	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.14	164	4842	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.89	188	9717	0.40	ng/ul	0.00
16) Chrysene-d12	21.09	240	7653	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	8017	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.86	152	604	0.04	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	1213	0.04	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.31	128	928	0.033	ng/ul#	82
5) 2-Methylnaphthalene	11.94	142	530	0.025	ng/ul	98
12) Phenanthrene	16.93	178	2091	0.062	ng/ul#	93
15) Fluoranthene	18.95	202	3251	0.084	ng/ul	94
17) Pyrene	19.32	202	3119	0.080	ng/ul#	94
18) Benzo(a)anthracene	21.07	228	1620	0.048	ng/ul#	82
19) Chrysene	21.12	228	2178	0.064	ng/ul#	85
21) Benzo(b)fluoranthene	22.58	252	3025m	0.081	ng/ul	
22) Benzo(k)fluoranthene	22.62	252	1207m	0.032	ng/ul	
23) Benzo(a)pyrene	23.12	252	1759	0.054	ng/ul#	32
24) Indeno(1,2,3-cd)pyrene	25.31	276	1597	0.045	ng/ul	91
26) Benzo(g,h,i)perylene	25.94	276	2032	0.068	ng/ul#	75

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

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