

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN051519\
 Data File : BN005696.D
 Acq On : 15 May 2019 15:24
 Operator : JU/SJ
 Sample : K2692-09
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A5118

Manual Integrations
 APPROVED

mohammad
 5/16/2019 11:38:29 AM

Quant Time: May 15 16:12:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN050719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 14 07:15:06 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	1069	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	4734	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	3510	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	9412	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	9657	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	9851	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	1857	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	7356	0.25	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.41	128	4882	0.390	ng/ul	95
5) 2-Methylnaphthalene	12.04	142	4001	0.468	ng/ul	99
7) Acenaphthylene	13.94	152	7514	0.425	ng/ul	99
8) Acenaphthene	14.28	153	6233	0.498	ng/ul	98
11) Pentachlorophenol	16.68	266	1746	0.716	ng/ul	88
13) Anthracene	17.12	178	9509	0.372	ng/ul	99
17) Pyrene	19.43	202	21044	0.590	ng/ul	96
19) Chrysene	21.26	228	16148	0.395	ng/ul	98
21) Benzo(b)fluoranthene	22.78	252	18429m	0.622	ng/ul	
22) Benzo(k)fluoranthene	22.83	252	30269	0.815	ng/ul	97
23) Benzo(a)pyrene	23.35	252	12563	0.380	ng/ul#	90
25) Dibenzo(a,h)anthracene	25.65	278	15071	0.492	ng/ul	97

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

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