

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022822\
 Data File : BN018755.D
 Acq On : 28 Feb 2022 12:04
 Operator : CG/JU
 Sample : N1656-13
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGQ20

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 03/01/2022
 Supervised By :Yogesh Patel 03/02/2022

Quant Time: Feb 28 23:54:38 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN021822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 18 13:39:55 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.913	152	5914	0.400	ng/ul	0.00
4) Naphthalene-d8	10.720	136	14235	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.546	164	7287	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.287	188	13523	0.400	ng/ul	0.00
17) Chrysene-d12	21.466	240	10209	0.400	ng/ul	# 0.00
23) Perylene-d12	23.863	264	9647	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.285	96	35441	5.277	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.309	152	6867	0.353	ng/ul	0.00
18) Fluoranthene-d10	19.311	212	15007	0.492	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.322	88	2222	0.323	ng/ul#	85
15) Phenanthrene	17.325	178	2650	0.060	ng/ul#	91
19) Fluoranthene	19.339	202	3693	0.086	ng/ul#	85
20) Pyrene	19.701	202	2563	0.058	ng/ul#	71
22) Chrysene	21.501	228	2511	0.061	ng/ul#	93
24) Benzo(b)fluoranthene	23.132	252	5583m	0.130	ng/ul	
25) Benzo(k)fluoranthene	23.176	252	1626m	0.038	ng/ul	
26) Benzo(a)pyrene	23.755	252	2108	0.056	ng/ul#	33
27) Indeno(1,2,3-cd)pyrene	26.344	276	3407	0.071	ng/ul#	88
28) Dibenzo(a,h)anthracene	26.358	278	813	0.021	ng/ul#	15
29) Benzo(g,h,i)perylene	27.112	276	3811	0.092	ng/ul#	84

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

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