

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042922\
 Data File : BN019588.D
 Acq On : 29 Apr 2022 12:54
 Operator : CG/JU
 Sample : N2469-01DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BFFF8DL

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 05/04/2022
 Supervised By :Yogesh Patel 05/05/2022

Quant Time: Apr 30 03:18:16 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN042922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 28 13:24:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.880	152	5575	0.400	ng/ul	0.00
4) Naphthalene-d8	10.667	136	16271	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.497	164	8573	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.234	188	16678	0.400	ng/ul	0.00
17) Chrysene-d12	21.416	240	14764	0.400	ng/ul	0.00
23) Perylene-d12	23.769	264	13737	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.349	96	3116	0.472	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	887	0.036	ng/ul	0.00
18) Fluoranthene-d10	19.257	212	1716	0.037	ng/ul	0.00
Target Compounds						
					Qvalue	
10) Acenaphthylene	14.215	152	3646	0.086	ng/ul	96
15) Phenanthrene	17.272	178	14680	0.239	ng/ul	99
16) Anthracene	17.365	178	4154	0.083	ng/ul	96
19) Fluoranthene	19.290	202	48951	0.724	ng/ul	99
20) Pyrene	19.652	202	52592	0.780	ng/ul	100
21) Benzo(a)anthracene	21.398	228	23367	0.423	ng/ul	96
22) Chrysene	21.451	228	36508	0.561	ng/ul	99
24) Benzo(b)fluoranthene	23.055	252	61548m	0.883	ng/ul	
25) Benzo(k)fluoranthene	23.093	252	20034m	0.298	ng/ul	
26) Benzo(a)pyrene	23.664	252	31452	0.530	ng/ul#	80
27) Indeno(1,2,3-cd)pyrene	26.183	276	28173	0.380	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.200	278	6262	0.101	ng/ul#	72
29) Benzo(g,h,i)perylene	26.927	276	25402	0.375	ng/ul	95

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

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