

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN12222\
 Data File : BN023354.D
 Acq On : 23 Dec 2022 00:18
 Operator : CG/JU
 Sample : N6003-18DL 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBXT8DL

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/23/2022
 Supervised By :mohammad ahmed 12/23/2022

Quant Time: Dec 23 01:26:02 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 22 02:28:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.998	152	6268	0.400	ng/ul	0.00
4) Naphthalene-d8	10.815	136	21493	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.645	164	13104	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.390	188	31236	0.400	ng/ul	0.00
17) Chrysene-d12	21.578	240	24514	0.400	ng/ul	# 0.00
23) Perylene-d12	24.028	264	19718	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.345	96	9157	1.345	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.399	152	3055	0.092	ng/ul	0.00
18) Fluoranthene-d10	19.420	212	7926	0.089	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.865	128	2163	0.034	ng/ul#	87
10) Acenaphthylene	14.368	152	3451	0.059	ng/ul#	93
15) Phenanthrene	17.432	178	9377	0.092	ng/ul	96
16) Anthracene	17.525	178	2457	0.029	ng/ul#	86
19) Fluoranthene	19.448	202	29972	0.252	ng/ul	98
20) Pyrene	19.811	202	23474m	0.196	ng/ul	
21) Benzo(a)anthracene	21.560	228	19913	0.209	ng/ul	95
22) Chrysene	21.616	228	22642	0.233	ng/ul#	92
24) Benzo(b)fluoranthene	23.279	252	31447m	0.308	ng/ul	
25) Benzo(k)fluoranthene	23.323	252	11144m	0.116	ng/ul	
26) Benzo(a)pyrene	23.916	252	16234	0.201	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	26.584	276	11891	0.126	ng/ul#	90
28) Dibenzo(a,h)anthracene	26.594	278	3139	0.043	ng/ul#	1
29) Benzo(g,h,i)perylene	27.372	276	8018	0.104	ng/ul#	85

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

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