

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082824\
 Data File : BN033643.D
 Acq On : 28 Aug 2024 13:32
 Operator : MA/JU
 Sample : P3711-10MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E24A0MSD

Quant Time: Aug 28 15:07:01 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN082724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 23:48:09 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.536	152	5594	0.400	ng/ul	0.00
4) Naphthalene-d8	10.299	136	15220	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.174	164	7554	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.927	188	15718	0.400	ng/ul	0.00
17) Chrysene-d12	21.138	240	12644	0.400	ng/ul	0.00
23) Perylene-d12	23.301	264	13871	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.148	96	1760	0.306	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.894	152	7069	0.324	ng/ul	0.00
18) Fluoranthene-d10	18.966	212	17050	0.452	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.177	88	5200	0.815	ng/ul#	87
5) Naphthalene	10.349	128	12326	0.303	ng/ul	100
7) 2-Methylnaphthalene	11.971	142	7615	0.295	ng/ul	100
8) 1-Methylnaphthalene	12.191	142	9010	0.340	ng/ul	99
10) Acenaphthylene	13.892	152	10665	0.311	ng/ul	99
11) Acenaphthene	14.239	153	8349	0.337	ng/ul	98
12) Fluorene	15.234	166	9442	0.341	ng/ul	100
14) Pentachlorophenol	16.581	266	4252	0.860	ng/ul	99
15) Phenanthrene	16.970	178	15732	0.339	ng/ul	100
16) Anthracene	17.058	178	13931	0.334	ng/ul	99
19) Fluoranthene	18.994	202	19050	0.382	ng/ul	99
20) Pyrene	19.361	202	20331	0.384	ng/ul	98
21) Benzo(a)anthracene	21.121	228	18977	0.381	ng/ul	99
22) Chrysene	21.176	228	19280	0.383	ng/ul	99
24) Benzo(b)fluoranthene	22.661	252	20009	0.372	ng/ul	99
25) Benzo(k)fluoranthene	22.702	252	20373	0.378	ng/ul	99
26) Benzo(a)pyrene	23.210	252	17726	0.410	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.448	276	24337	0.375	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.465	278	19021	0.377	ng/ul	100
29) Benzo(g,h,i)perylene	26.102	276	19130	0.377	ng/ul	99

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

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