

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090324\
 Data File : BN033743.D
 Acq On : 03 Sep 2024 16:21
 Operator : MA/JU
 Sample : P3809-09MS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 YE3E3MS

Quant Time: Sep 04 01:41:46 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 : Wed Sep 04 01:39:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.523	152	5981	0.400	ng/ul	0.00
4) Naphthalene-d8	10.283	136	16115	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.160	164	8071	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.910	188	17212	0.400	ng/ul	0.00
17) Chrysene-d12	21.121	240	14779	0.400	ng/ul	0.00
23) Perylene-d12	23.278	264	15918	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.135	96	1954	0.318	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.878	152	8741	0.378	ng/ul	0.00
18) Fluoranthene-d10	18.952	212	22372	0.508	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.164	88	64098	9.395	ng/ul#	75
5) Naphthalene	10.327	128	14686	0.341	ng/ul	100
7) 2-Methylnaphthalene	11.955	142	9372	0.343	ng/ul	99
8) 1-Methylnaphthalene	12.175	142	11327	0.403	ng/ul	99
10) Acenaphthylene	13.874	152	13447	0.367	ng/ul	99
11) Acenaphthene	14.225	153	10471	0.396	ng/ul	98
12) Fluorene	15.215	166	12354	0.418	ng/ul	100
14) Pentachlorophenol	16.568	266	4649	0.858	ng/ul	99
15) Phenanthrene	16.953	178	21187	0.417	ng/ul	99
16) Anthracene	17.045	178	19347	0.424	ng/ul	99
19) Fluoranthene	18.980	202	24605	0.422	ng/ul	98
20) Pyrene	19.343	202	26163	0.423	ng/ul	99
21) Benzo(a)anthracene	21.106	228	24612	0.423	ng/ul	98
22) Chrysene	21.156	228	24298	0.413	ng/ul	99
24) Benzo(b)fluoranthene	22.634	252	25705	0.416	ng/ul	99
25) Benzo(k)fluoranthene	22.678	252	25971	0.420	ng/ul	100
26) Benzo(a)pyrene	23.181	252	22396	0.451	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.415	276	29517	0.396	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.428	278	22899	0.396	ng/ul	100
29) Benzo(g,h,i)perylene	26.062	276	23084	0.396	ng/ul	99

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

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