

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071224\
 Data File : BN032660.D
 Acq On : 12 Jul 2024 06:24
 Operator : MA/JU
 Sample : P3138-03MS
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 YDZF8MS

Quant Time: Jul 12 06:48:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN071224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 01:40:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.557	152	5548	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.327	136	17631	0.400	ng/ul	-0.03
9) Acenaphthene-d10	14.202	164	9843	0.400	ng/ul	-0.01
13) Phenanthrene-d10	16.961	188	16543	0.400	ng/ul	-0.03
17) Chrysene-d12	21.170	240	13125	0.400	ng/ul	-0.02
23) Perylene-d12	23.365	264	15487	0.400	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.105	96	1547	0.232	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.927	152	9421	0.354	ng/ul	-0.04
18) Fluoranthene-d10	18.999	212	17935	0.419	ng/ul	-0.03
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.135	88	5845	0.770	ng/ul#	83
5) Naphthalene	10.376	128	17809	0.376	ng/ul	98
7) 2-Methylnaphthalene	11.999	142	11366	0.385	ng/ul	96
8) 1-Methylnaphthalene	12.218	142	11372	0.372	ng/ul	100
10) Acenaphthylene	13.924	152	15486	0.389	ng/ul	100
11) Acenaphthene	14.267	153	11897	0.362	ng/ul	99
12) Fluorene	15.261	166	12890	0.343	ng/ul	99
14) Pentachlorophenol	16.627	266	4618	1.591	ng/ul	99
15) Phenanthrene	17.003	178	19739	0.438	ng/ul	99
16) Anthracene	17.096	178	18582	0.505	ng/ul	99
19) Fluoranthene	19.031	202	22362	0.419	ng/ul	99
20) Pyrene	19.398	202	23675	0.429	ng/ul	99
21) Benzo(a)anthracene	21.156	228	20876	0.500	ng/ul	98
22) Chrysene	21.208	228	22559	0.373	ng/ul	97
24) Benzo(b)fluoranthene	22.707	252	22626	0.414	ng/ul	95
25) Benzo(k)fluoranthene	22.751	252	22880	0.344	ng/ul	95
26) Benzo(a)pyrene	23.272	252	20913	0.442	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.579	276	28126	0.439	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.582	278	21990	0.443	ng/ul	96
29) Benzo(g,h,i)perylene	26.253	276	22127	0.413	ng/ul	99

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

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