

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034950.D
 Acq On : 09 Nov 2024 20:37
 Operator : RC/JU
 Sample : P4744-18MS
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GCP72MS

Quant Time: Nov 09 22:53:43 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 09 00:09:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.571	152	3681	0.400	ng/ul	0.00
4) Naphthalene-d8	10.334	136	10170	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.203	164	8727	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.950	188	18838	0.400	ng/ul #	0.00
17) Chrysene-d12	21.145	240	16238	0.400	ng/ul #	0.00
23) Perylene-d12	23.311	264	14021	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	776	0.222	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.928	152	6105	0.392	ng/ul	0.00
18) Fluoranthene-d10	18.981	212	21794	0.523	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.183	88	2493	0.649	ng/ul#	85
5) Naphthalene	10.383	128	9740	0.383	ng/ul	100
7) 2-Methylnaphthalene	12.005	142	7828	0.412	ng/ul	98
8) 1-Methylnaphthalene	12.225	142	7764	0.407	ng/ul	98
10) Acenaphthylene	13.921	152	12309	0.345	ng/ul	97
11) Acenaphthene	14.268	153	9085	0.336	ng/ul	98
12) Fluorene	15.258	166	13438	0.391	ng/ul	97
14) Pentachlorophenol	16.603	266	6510	1.495	ng/ul	97
15) Phenanthrene	16.992	178	20361	0.404	ng/ul#	93
16) Anthracene	17.080	178	19460	0.411	ng/ul	93
19) Fluoranthene	19.014	202	25982	0.479	ng/ul#	94
20) Pyrene	19.376	202	28590	0.508	ng/ul#	94
21) Benzo(a)anthracene	21.131	228	24355	0.414	ng/ul#	93
22) Chrysene	21.180	228	23038	0.384	ng/ul#	92
24) Benzo(b)fluoranthene	22.668	252	20356	0.413	ng/ul	91
25) Benzo(k)fluoranthene	22.709	252	20348	0.394	ng/ul#	91
26) Benzo(a)pyrene	23.217	252	17535	0.398	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	25.467	276	20152	0.374	ng/ul#	94
28) Dibenzo(a,h)anthracene	25.480	278	15871	0.378	ng/ul#	94
29) Benzo(g,h,i)perylene	26.117	276	15451	0.375	ng/ul#	97

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

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