

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070723\
 Data File : BN026357.D
 Acq On : 07 Jul 2023 13:53
 Operator : MA/JU
 Sample : 03284-11MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DCGZ6MSD

Quant Time: Jul 07 15:54:35 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 07 14:46:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.938	152	4784	0.400	ng/ul	0.00
4) Naphthalene-d8	10.746	136	16834	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.579	164	9193	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.329	188	14182	0.400	ng/ul	0.00
17) Chrysene-d12	21.528	240	9422	0.400	ng/ul	0.00
23) Perylene-d12	23.995	264	10332	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	8187	1.520	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.336	152	7478	0.289	ng/ul	0.00
18) Fluoranthene-d10	19.357	212	13462	0.375	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.343	88	5610	0.981	ng/ul	93
5) Naphthalene	10.796	128	16018	0.332	ng/ul	99
7) 2-Methylnaphthalene	12.407	142	9554	0.311	ng/ul	99
8) 1-Methylnaphthalene	12.622	142	10251	0.322	ng/ul	100
10) Acenaphthylene	14.301	152	14148	0.329	ng/ul	99
11) Acenaphthene	14.639	153	11215	0.318	ng/ul	99
12) Fluorene	15.629	166	12026	0.298	ng/ul	99
14) Pentachlorophenol	16.991	266	2216	0.436	ng/ul	99
15) Phenanthrene	17.371	178	16254	0.347	ng/ul	99
16) Anthracene	17.464	178	14719	0.354	ng/ul	97
19) Fluoranthene	19.390	202	18626	0.370	ng/ul	98
20) Pyrene	19.752	202	19725	0.377	ng/ul	97
21) Benzo(a)anthracene	21.513	228	13719	0.353	ng/ul	99
22) Chrysene	21.566	228	14796	0.366	ng/ul	99
24) Benzo(b)fluoranthene	23.240	252	14592	0.331	ng/ul	92
25) Benzo(k)fluoranthene	23.287	252	15825	0.357	ng/ul	94
26) Benzo(a)pyrene	23.889	252	13045	0.341	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	26.580	276	17318	0.414	ng/ul#	44
28) Dibenzo(a,h)anthracene	26.586	278	13556	0.428	ng/ul	99
29) Benzo(g,h,i)perylene	27.365	276	15787	0.424	ng/ul	98

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

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