

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080224\
 Data File : BN033179.D
 Acq On : 01 Aug 2024 22:30
 Operator : MA/JU
 Sample : P3401-05MSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E1GG3MSD

Quant Time: Aug 01 23:34:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN072524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 30 14:49:49 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.512	152	2733	0.400	ng/ul	-0.04
4) Naphthalene-d8	10.267	136	9683	0.400	ng/ul	-0.05
9) Acenaphthene-d10	14.134	164	5807	0.400	ng/ul	-0.03
13) Phenanthrene-d10	16.911	188	12969	0.400	ng/ul	-0.03
17) Chrysene-d12	21.131	240	10269	0.400	ng/ul	-0.02
23) Perylene-d12	23.323	264	9992	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.052	96	923	0.267	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.857	152	4256	0.269	ng/ul	-0.07
18) Fluoranthene-d10	18.953	212	12407	0.427	ng/ul	-0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	3.086	88	3238	0.822	ng/ul#	73
5) Naphthalene	10.317	128	10014	0.403	ng/ul	96
7) 2-Methylnaphthalene	11.934	142	6387	0.379	ng/ul	98
8) 1-Methylnaphthalene	12.148	142	5416	0.297	ng/ul	98
10) Acenaphthylene	13.866	152	9010	0.379	ng/ul	100
11) Acenaphthene	14.199	153	7800	0.436	ng/ul	99
12) Fluorene	15.212	166	8893	0.427	ng/ul	99
14) Pentachlorophenol	16.595	266	2078	0.787	ng/ul	99
15) Phenanthrene	16.954	178	14501	0.417	ng/ul	100
16) Anthracene	17.055	178	12444	0.398	ng/ul	99
19) Fluoranthene	18.986	202	18101	0.535	ng/ul	99
20) Pyrene	19.353	202	19029	0.525	ng/ul	99
21) Benzo(a)anthracene	21.119	228	14116	0.403	ng/ul	99
22) Chrysene	21.169	228	17182	0.420	ng/ul	99
24) Benzo(b)fluoranthene	22.668	252	12420	0.426	ng/ul	84
25) Benzo(k)fluoranthene	22.712	252	15964	0.456	ng/ul#	87
26) Benzo(a)pyrene	23.238	252	10615	0.386	ng/ul#	78
27) Indeno(1,2,3-cd)pyrene	25.534	276	14612	0.350	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.524	278	11671	0.353	ng/ul	95
29) Benzo(g,h,i)perylene	26.214	276	10725	0.323	ng/ul	97

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

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