

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090322\
 Data File : BM036493.D
 Acq On : 02 Sep 2022 16:13
 Operator : CG/JU
 Sample : PB147415BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS415

Quant Time: Sep 02 16:52:45 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM083022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 29 17:45:20 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.737	152	3744	0.400	ng/ul	0.00
4) Naphthalene-d8	10.534	136	15040	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.381	164	8396	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.127	188	17073	0.400	ng/ul	0.00
17) Chrysene-d12	21.316	240	15189	0.400	ng/ul	# 0.00
23) Perylene-d12	23.607	264	11984	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.139	96	3226	0.593	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.134	152	7495	0.319	ng/ul	0.00
18) Fluoranthene-d10	19.159	212	15529	0.329	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.172	88	8316	1.609	ng/ul#	81
5) Naphthalene	10.583	128	12879	0.313	ng/ul	97
7) 2-Methylnaphthalene	12.206	142	7965	0.316	ng/ul	95
8) 1-Methylnaphthalene	12.420	142	8416	0.328	ng/ul	99
10) Acenaphthylene	14.103	152	11101	0.348	ng/ul	99
11) Acenaphthene	14.441	153	8871	0.315	ng/ul	99
12) Fluorene	15.435	166	9786	0.312	ng/ul	100
14) Pentachlorophenol	16.794	266	1845	0.600	ng/ul	97
15) Phenanthrene	17.170	178	16083	0.323	ng/ul	100
16) Anthracene	17.267	178	13620	0.325	ng/ul	99
19) Fluoranthene	19.191	202	18623	0.341	ng/ul	98
20) Pyrene	19.554	202	19352	0.343	ng/ul	97
21) Benzo(a)anthracene	21.301	228	15715	0.342	ng/ul	99
22) Chrysene	21.354	228	19195	0.374	ng/ul	99
24) Benzo(b)fluoranthene	22.911	252	17528	0.368	ng/ul	93
25) Benzo(k)fluoranthene	22.955	252	16919	0.355	ng/ul#	91
26) Benzo(a)pyrene	23.511	252	16111	0.413	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	25.977	276	17506	0.377	ng/ul#	93
28) Dibenzo(a,h)anthracene	26.001	278	14395	0.362	ng/ul#	90
29) Benzo(g,h,i)perylene	26.702	276	15840	0.358	ng/ul	95

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

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