

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080921\
 Data File : BM031650.D
 Acq On : 10 Aug 2021 20:31
 Operator : CG/JU
 Sample : M3169-05DL 5X
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C00A4DL

Manual Integrations
 APPROVED

mohammad
 8/11/2021 9:25:43 PM

Quant Time: Aug 11 05:58:21 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 10 18:36:58 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.576	152	1505	0.400	ng/ul	0.00
4) Naphthalene-d8	10.338	136	5591	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.209	164	3146	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.961	188	5940	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.156	240	4655	0.400	ng/ul	# 0.00
23) Perylene-d12	23.314	264	4802	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.189	96	1413	0.845	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.930	152	484	0.051	ng/ul	0.00
18) Fluoranthene-d10	18.993	212	1043	0.067	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.388	128	6987	0.420	ng/ul	99
7) 2-Methylnaphthalene	12.007	142	3958	0.348	ng/ul	99
8) 1-Methylnaphthalene	12.232	142	2452	0.218	ng/ul	98
10) Acenaphthylene	13.932	152	11101	0.828	ng/ul	99
11) Acenaphthene	14.273	153	1838	0.163	ng/ul	98
12) Fluorene	15.266	166	4325	0.333	ng/ul#	96
15) Phenanthrene	17.002	178	86193	4.570	ng/ul	97
16) Anthracene	17.093	178	19759	1.119	ng/ul	99
19) Fluoranthene	19.024	202	163595	8.230	ng/ul	97
20) Pyrene	19.386	202	135988	6.745	ng/ul	95
21) Benzo(a)anthracene	21.139	228	66950	3.562	ng/ul	97
22) Chrysene	21.190	228	66169	3.402	ng/ul	96
24) Benzo(b)fluoranthene	22.674	252	86613m	4.003	ng/ul	
25) Benzo(k)fluoranthene	22.713	252	31354m	1.403	ng/ul	
26) Benzo(a)pyrene	23.220	252	62551	3.301	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	25.446	276	42181	1.722	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.451	278	10409	0.527	ng/ul#	85
29) Benzo(g,h,i)perylene	26.096	276	35846	1.716	ng/ul	98

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

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