

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM103124\
 Data File : BM048379.D
 Acq On : 01 Nov 2024 04:41
 Operator : RC/JU
 Sample : P4592-02
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F7H02

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/04/2024
 Supervised By :mohammad ahmed 11/04/2024

Quant Time: Nov 03 07:05:23 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM103124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 31 14:58:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.708	152	4893	0.400	ng/ul	-0.04
4) Naphthalene-d8	10.495	136	15750	0.400	ng/ul	#-0.03
9) Acenaphthene-d10	14.348	164	8608	0.400	ng/ul	-0.02
13) Phenanthrene-d10	17.094	188	17875m	0.400	ng/ul	-0.03
17) Chrysene-d12	21.289	240	15050m	0.400	ng/ul	0.00
23) Perylene-d12	23.516	264	12706m	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	21184	3.525	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.118	152	5538	0.276	ng/ul	-0.02
18) Fluoranthene-d10	19.126	212	12892	0.337	ng/ul	-0.01
Target Compounds						
					Qvalue	
11) Acenaphthene	14.413	153	1953	0.073	ng/ul	99
12) Fluorene	15.403	166	2529	0.085	ng/ul	98
15) Phenanthrene	17.136	178	5557m	0.125	ng/ul	
16) Anthracene	17.224	178	1173m	0.028	ng/ul	
19) Fluoranthene	19.154	202	7199	0.139	ng/ul	95
20) Pyrene	19.512	202	7227m	0.130	ng/ul	
21) Benzo(a)anthracene	21.278	228	3249	0.072	ng/ul#	89
22) Chrysene	21.321	228	3607	0.058	ng/ul#	91
24) Benzo(b)fluoranthene	22.841	252	5133m	0.118	ng/ul	
25) Benzo(k)fluoranthene	22.876	252	2150m	0.042	ng/ul	
26) Benzo(a)pyrene	23.408	252	2395	0.059	ng/ul#	48
27) Indeno(1,2,3-cd)pyrene	25.826	276	1908	0.033	ng/ul#	83
29) Benzo(g,h,i)perylene	26.520	276	1826	0.038	ng/ul#	64

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

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