

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122024\
 Data File : BM049107.D
 Acq On : 20 Dec 2024 19:13
 Operator : RC/JU
 Sample : P5265-02DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E28Y7DL

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/23/2024
 Supervised By :mohammad ahmed 12/24/2024

Quant Time: Dec 20 22:10:54 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM121824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 18 15:33:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.560	152	7744	0.400	ng/ul	0.00
4) Naphthalene-d8	10.321	136	22912	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.197	164	13076	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.943	188	26517	0.400	ng/ul	0.00
17) Chrysene-d12	21.146	240	22815	0.400	ng/ul	0.00
23) Perylene-d12	23.303	264	25151	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.151	96	4838	0.542	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.926	152	1115	0.036	ng/ul	0.00
18) Fluoranthene-d10	18.979	212	2186	0.034	ng/ul	0.00
Target Compounds						
11) Acenaphthene	14.257	153	1115	0.028	ng/ul	Qvalue 94
12) Fluorene	15.252	166	897	0.020	ng/ul#	91
15) Phenanthrene	16.985	178	14943	0.216	ng/ul	99
16) Anthracene	17.078	178	3162	0.050	ng/ul#	90
19) Fluoranthene	19.011	202	28801	0.349	ng/ul	100
20) Pyrene	19.374	202	21315	0.242	ng/ul	100
21) Benzo(a)anthracene	21.128	228	12220	0.154	ng/ul	97
22) Chrysene	21.181	228	13017	0.153	ng/ul	95
24) Benzo(b)fluoranthene	22.665	252	16193m	0.202	ng/ul	
25) Benzo(k)fluoranthene	22.703	252	7083m	0.080	ng/ul	
26) Benzo(a)pyrene	23.209	252	6737	0.087	ng/ul#	63
27) Indeno(1,2,3-cd)pyrene	25.443	276	6320	0.058	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.457	278	1957	0.023	ng/ul#	11

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

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