

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042424\
 Data File : BM045531.D
 Acq On : 24 Apr 2024 16:56
 Operator : MA/JU
 Sample : P2164-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COSA0

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 04/25/2024
 Supervised By :mohammad ahmed 04/27/2024

Quant Time: Apr 24 17:36:09 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM042024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 24 05:17:39 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.551	152	1003	0.400	ng/ul	-0.01	
4) Naphthalene-d8	10.314	136	2971	0.400	ng/ul	#-0.02	
9) Acenaphthene-d10	14.200	164	1649	0.400	ng/ul	0.00	
13) Phenanthrene-d10	16.963	188	3556m	0.400	ng/ul	-0.02	
17) Chrysene-d12	21.199	240	3225	0.400	ng/ul	# 0.00	
23) Perylene-d12	23.388	264	3549m	0.400	ng/ul	0.00	
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.163	96	4779	3.555	ng/ul	-0.01	
6) 2-Methylnaphthalene-d10	11.936	152	972	0.248	ng/ul	0.00	
18) Fluoranthene-d10	19.010	212	2141	0.211	ng/ul	0.00	
Target Compounds							
10) Acenaphthylene	13.923	152	254	0.031	ng/ul#		47
15) Phenanthrene	17.005	178	989	0.095	ng/ul		94
19) Fluoranthene	19.038	202	2013	0.135	ng/ul		99
20) Pyrene	19.400	202	2318m	0.152	ng/ul		
21) Benzo(a)anthracene	21.190	228	812	0.079	ng/ul#		89
22) Chrysene	21.237	228	1148	0.072	ng/ul#		90
24) Benzo(b)fluoranthene	22.727	252	1564m	0.131	ng/ul		
25) Benzo(k)fluoranthene	22.762	252	685m	0.045	ng/ul		
26) Benzo(a)pyrene	23.279	252	1078	0.087	ng/ul#		52
27) Indeno(1,2,3-cd)pyrene	25.608	276	1024	0.057	ng/ul#		85
29) Benzo(g,h,i)perylene	26.262	276	1070	0.066	ng/ul#		74

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

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