

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070323\
 Data File : BM040587.D
 Acq On : 03 Jul 2023 22:32
 Operator : MA/JU
 Sample : 03242-06DL 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCGA4DL

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/04/2023
 Supervised By :Sohil Jodhani 07/04/2023

Quant Time: Jul 04 01:30:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 03 13:40:38 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.923	152	6014	0.400	ng/ul	0.00
4) Naphthalene-d8	10.728	136	22726	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.568	164	11118	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.315	188	21798	0.400	ng/ul	0.00
17) Chrysene-d12	21.497	240	12963	0.400	ng/ul	0.00
23) Perylene-d12	23.906	264	12209	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.321	96	6391	0.677	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.323	152	1528	0.048	ng/ul	0.00
18) Fluoranthene-d10	19.342	212	2309	0.056	ng/ul	0.00
Target Compounds						
					Qvalue	
10) Acenaphthylene	14.286	152	3705	0.075	ng/ul#	91
15) Phenanthrene	17.357	178	8471	0.125	ng/ul	97
16) Anthracene	17.446	178	3446	0.061	ng/ul#	86
19) Fluoranthene	19.370	202	38785	0.704	ng/ul	98
20) Pyrene	19.737	202	30660	0.519	ng/ul	98
21) Benzo(a)anthracene	21.483	228	23229	0.530	ng/ul	96
22) Chrysene	21.535	228	22253	0.445	ng/ul	97
24) Benzo(b)fluoranthene	23.175	252	34164	0.723	ng/ul	98
25) Benzo(k)fluoranthene	23.210	252	11676m	0.251	ng/ul	
26) Benzo(a)pyrene	23.797	252	19177	0.451	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.407	276	15115	0.255	ng/ul#	88
28) Dibenzo(a,h)anthracene	26.417	278	4420	0.095	ng/ul#	60
29) Benzo(g,h,i)perylene	27.178	276	13852	0.267	ng/ul	94

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

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