

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010323\
 Data File : BM038357.D
 Acq On : 03 Jan 2023 12:41
 Operator : CG/JU
 Sample : N5388-02
 Misc : SFAM-SIM-MDL-W-02
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-WATER-QT4-2022-08

Quant Time: Jan 03 13:12:00 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM123022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 30 01:43:54 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 01/04/2023
 Supervised By :mohammad ahmed 01/05/2023

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.009	152	1365	0.400	ng/ul	0.00
4) Naphthalene-d8	10.823	136	4101	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.638	164	2162	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.379	188	4408	0.400	ng/ul	0.00
17) Chrysene-d12	21.550	240	4404	0.400	ng/ul	0.00
23) Perylene-d12	23.979	264	5601	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	1239	0.609	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.407	152	2041	0.332	ng/ul	0.00
18) Fluoranthene-d10	19.399	212	4879	0.391	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.410	88	177m	0.085	ng/ul	
5) Naphthalene	10.872	128	672	0.061	ng/ul#	71
7) 2-Methylnaphthalene	12.478	142	388	0.058	ng/ul	99
8) 1-Methylnaphthalene	12.693	142	397	0.058	ng/ul	96
10) Acenaphthylene	14.365	152	667	0.066	ng/ul#	81
11) Acenaphthene	14.698	153	484	0.065	ng/ul	94
12) Fluorene	15.688	166	521	0.062	ng/ul#	90
14) Pentachlorophenol	17.054	266	96	0.061	ng/ul	97
15) Phenanthrene	17.426	178	821	0.063	ng/ul#	82
16) Anthracene	17.518	178	712	0.059	ng/ul#	77
19) Fluoranthene	19.431	202	1039	0.068	ng/ul#	80
20) Pyrene	19.794	202	1100	0.068	ng/ul#	81
21) Benzo(a)anthracene	21.536	228	945	0.064	ng/ul#	88
22) Chrysene	21.588	228	1011	0.065	ng/ul#	91
24) Benzo(b)fluoranthene	23.237	252	1102	0.054	ng/ul#	45
25) Benzo(k)fluoranthene	23.284	252	1161	0.057	ng/ul#	46
26) Benzo(a)pyrene	23.874	252	1045	0.054	ng/ul#	34
27) Indeno(1,2,3-cd)pyrene	26.525	276	1703	0.061	ng/ul#	77
28) Dibenzo(a,h)anthracene	26.538	278	1367	0.062	ng/ul#	34
29) Benzo(g,h,i)perylene	27.310	276	1575	0.063	ng/ul#	80

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

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