

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080221\
 Data File : BM031475.D
 Acq On : 05 Aug 2021 08:39
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4095

Quant Time: Aug 05 09:12:06 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 05 07:10:43 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.594	152	1443	0.400	ng/ul	0.00
4) Naphthalene-d8	10.352	136	5305	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.224	164	3018	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.971	188	5753	0.400	ng/ul	0.00
17) Chrysene-d12	21.166	240	4482	0.400	ng/ul	0.00
23) Perylene-d12	23.324	264	4227	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.196	96	568	0.354	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.948	152	3112	0.348	ng/ul	0.00
18) Fluoranthene-d10	19.005	212	5381	0.360	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.227	88	602	0.372	ng/ul#	87
5) Naphthalene	10.401	128	5663	0.358	ng/ul	99
7) 2-Methylnaphthalene	12.020	142	3735	0.346	ng/ul	99
8) 1-Methylnaphthalene	12.245	142	3679	0.345	ng/ul	100
10) Acenaphthylene	13.941	152	4581	0.356	ng/ul	99
11) Acenaphthene	14.284	153	3751	0.347	ng/ul	98
12) Fluorene	15.282	166	4115	0.331	ng/ul	99
14) Pentachlorophenol	16.631	266	599	0.325	ng/ul	99
15) Phenanthrene	17.013	178	6277	0.344	ng/ul	100
16) Anthracene	17.107	178	5810	0.340	ng/ul	99
19) Fluoranthene	19.035	202	6880	0.359	ng/ul	97
20) Pyrene	19.397	202	7050	0.363	ng/ul	97
21) Benzo(a)anthracene	21.152	228	6222	0.344	ng/ul	99
22) Chrysene	21.202	228	6498	0.347	ng/ul	98
24) Benzo(b)fluoranthene	22.682	252	6509	0.342	ng/ul	96
25) Benzo(k)fluoranthene	22.725	252	6608	0.336	ng/ul	98
26) Benzo(a)pyrene	23.232	252	5756	0.345	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	25.463	276	7174	0.333	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.478	278	5712	0.329	ng/ul	99
29) Benzo(g,h,i)perylene	26.115	276	6170	0.335	ng/ul	98

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

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