

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062821\
 Data File : BM030693.D
 Acq On : 29 Jun 2021 16:24
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4714

Quant Time: Jun 29 17:01:42 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 28 14:33:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.791	152	2195	0.400	ng/ul	0.00
4) Naphthalene-d8	10.576	136	8482	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.417	164	4756	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.155	188	9025	0.400	ng/ul	0.00
17) Chrysene-d12	21.328	240	8486	0.400	ng/ul	0.00
23) Perylene-d12	23.585	264	8997	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.286	96	878	0.367	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.164	152	5122	0.384	ng/ul	0.00
18) Fluoranthene-d10	19.176	212	9679	0.398	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.317	88	920	0.372	ng/ul#	84
5) Naphthalene	10.621	128	9040	0.374	ng/ul	99
7) 2-Methylnaphthalene	12.241	142	6238	0.383	ng/ul	99
8) 1-Methylnaphthalene	12.456	142	6071	0.383	ng/ul	100
10) Acenaphthylene	14.136	152	8967	0.422	ng/ul	99
11) Acenaphthene	14.478	153	6415	0.373	ng/ul	99
12) Fluorene	15.467	166	7052	0.372	ng/ul	99
14) Pentachlorophenol	16.822	266	977	0.332	ng/ul	95
15) Phenanthrene	17.197	178	12483	0.411	ng/ul	99
16) Anthracene	17.290	178	10808	0.407	ng/ul	99
19) Fluoranthene	19.207	202	17931	0.479	ng/ul	99
20) Pyrene	19.569	202	17747	0.471	ng/ul	98
21) Benzo(a)anthracene	21.314	228	14290	0.422	ng/ul	96
22) Chrysene	21.364	228	14996	0.396	ng/ul	100
24) Benzo(b)fluoranthene	22.904	252	15282	0.383	ng/ul	96
25) Benzo(k)fluoranthene	22.948	252	14318	0.348	ng/ul	97
26) Benzo(a)pyrene	23.486	252	13167	0.388	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	25.865	276	16640	0.377	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.875	278	13210	0.373	ng/ul	97
29) Benzo(g,h,i)perylene	26.563	276	14856	0.387	ng/ul	98

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

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