

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090121\
 Data File : BN016123.D
 Acq On : 01 Sep 2021 10:14
 Operator : CG/JU
 Sample : SST0.818
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.8218

Quant Time: Sep 01 10:44:23 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 01 10:18:28 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	3116	0.400	ng/ul	0.00
4) Naphthalene-d8	10.661	136	12068	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.501	164	6634	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.246	188	13480	0.400	ng/ul	0.00
17) Chrysene-d12	21.433	240	13303	0.400	ng/ul	0.00
23) Perylene-d12	23.824	264	11750	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	2895	0.681	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.250	152	13423	0.729	ng/ul	0.00
18) Fluoranthene-d10	19.276	212	27552	0.778	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	3243	0.736	ng/ul	92
5) Naphthalene	10.710	128	25037	0.810	ng/ul	99
7) 2-Methylnaphthalene	12.322	142	16569	0.789	ng/ul	100
8) 1-Methylnaphthalene	12.542	142	16482	0.803	ng/ul	100
10) Acenaphthylene	14.219	152	19594	0.801	ng/ul	99
11) Acenaphthene	14.561	153	16382	0.823	ng/ul	99
12) Fluorene	15.547	166	18851	0.814	ng/ul	99
14) Pentachlorophenol	16.896	266	1649	0.453	ng/ul	100
15) Phenanthrene	17.288	178	30281	0.816	ng/ul	99
16) Anthracene	17.377	178	26440	0.807	ng/ul	98
19) Fluoranthene	19.308	202	34645	0.851	ng/ul	99
20) Pyrene	19.671	202	34510	0.839	ng/ul	97
21) Benzo(a)anthracene	21.418	228	31419	0.749	ng/ul	100
22) Chrysene	21.471	228	35459	0.810	ng/ul	99
24) Benzo(b)fluoranthene	23.093	252	33941	0.806	ng/ul	96
25) Benzo(k)fluoranthene	23.140	252	35016	0.839	ng/ul	97
26) Benzo(a)pyrene	23.713	252	29127	0.794	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.290	276	38510	0.800	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.313	278	31929	0.793	ng/ul	98
29) Benzo(g,h,i)perylene	27.051	276	33266	0.813	ng/ul	98

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

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