

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102721\
 Data File : BN017134.D
 Acq On : 27 Oct 2021 19:32
 Operator : CG/JU
 Sample : SST0.832
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.832

Quant Time: Oct 28 01:06:42 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 28 01:03:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.531	152	4894	0.400	ng/ul	0.00
4) Naphthalene-d8	10.300	136	18859	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.185	164	11733	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.936	188	26054	0.400	ng/ul	0.00
17) Chrysene-d12	21.148	240	27116	0.400	ng/ul	0.00
23) Perylene-d12	23.354	264	26713	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.059	96	4183	0.787	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.906	152	21400	0.796	ng/ul	0.00
18) Fluoranthene-d10	18.976	212	54516	0.814	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.093	88	4509	0.704	ng/ul#	91
5) Naphthalene	10.350	128	40320	0.706	ng/ul	99
7) 2-Methylnaphthalene	11.978	142	28056	0.789	ng/ul	100
8) 1-Methylnaphthalene	12.203	142	28394	0.785	ng/ul	100
10) Acenaphthylene	13.898	152	36185	0.763	ng/ul	99
11) Acenaphthene	14.245	153	31597	0.776	ng/ul	100
12) Fluorene	15.240	166	37049	0.787	ng/ul	99
14) Pentachlorophenol	16.594	266	4308	0.686	ng/ul	100
15) Phenanthrene	16.978	178	63400	0.766	ng/ul	99
16) Anthracene	17.071	178	53701	0.769	ng/ul	99
19) Fluoranthene	19.009	202	75077	0.808	ng/ul	98
20) Pyrene	19.371	202	75473	0.790	ng/ul	97
21) Benzo(a)anthracene	21.133	228	68096	0.770	ng/ul	99
22) Chrysene	21.186	228	77007	0.766	ng/ul	100
24) Benzo(b)fluoranthene	22.691	252	77921	0.775	ng/ul	98
25) Benzo(k)fluoranthene	22.737	252	87518	0.790	ng/ul	99
26) Benzo(a)pyrene	23.255	252	68569	0.763	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.570	276	98837	0.774	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.590	278	78797	0.777	ng/ul	99
29) Benzo(g,h,i)perylene	26.244	276	82488	0.746	ng/ul	99

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

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