

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122122\
 Data File : BN023309.D
 Acq On : 21 Dec 2022 11:23
 Operator : CG/JU
 Sample : SST0.861
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.8219

Quant Time: Dec 21 12:00:05 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 21 11:52:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.998	152	6532	0.400	ng/u1	0.00
4) Naphthalene-d8	10.815	136	20102	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.645	164	11665	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.390	188	24644	0.400	ng/u1	0.00
17) Chrysene-d12	21.577	240	18935	0.400	ng/u1	0.00
23) Perylene-d12	24.026	264	12772	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.349	96	5779	0.795	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.404	152	26274	0.854	ng/u1	0.00
18) Fluoranthene-d10	19.420	212	56657	0.825	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.382	88	5967	0.807	ng/u1	93
5) Naphthalene	10.865	128	49676	0.845	ng/u1	100
7) 2-Methylnaphthalene	12.476	142	32062	0.860	ng/u1	100
8) 1-Methylnaphthalene	12.696	142	32736	0.847	ng/u1	100
10) Acenaphthylene	14.368	152	43819	0.868	ng/u1	99
11) Acenaphthene	14.710	153	36738	0.849	ng/u1	99
12) Fluorene	15.695	166	41885	0.873	ng/u1	98
14) Pentachlorophenol	17.040	266	5378	0.877	ng/u1	99
15) Phenanthrene	17.432	178	67831	0.855	ng/u1	100
16) Anthracene	17.525	178	57542	0.900	ng/u1	99
19) Fluoranthene	19.448	202	76250	0.823	ng/u1	100
20) Pyrene	19.810	202	75435	0.797	ng/u1	100
21) Benzo(a)anthracene	21.562	228	59324	0.793	ng/u1	100
22) Chrysene	21.615	228	61208	0.799	ng/u1	100
24) Benzo(b)fluoranthene	23.275	252	55591	0.844	ng/u1	97
25) Benzo(k)fluoranthene	23.322	252	53199	0.874	ng/u1	96
26) Benzo(a)pyrene	23.912	252	43568	0.833	ng/u1	95
27) Indeno(1,2,3-cd)pyrene	26.579	276	50373	0.825	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.595	278	39245	0.832	ng/u1	98
29) Benzo(g,h,i)perylene	27.367	276	40529	0.807	ng/u1	99

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

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