

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090121\
 Data File : BN016124.D
 Acq On : 01 Sep 2021 10:50
 Operator : CG/JU
 Sample : SSTD1.619
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.6219

Quant Time: Sep 01 11:20:58 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	2855	0.400	ng/ul	0.00
4) Naphthalene-d8	10.661	136	11234	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.497	164	6391	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.246	188	13125	0.400	ng/ul	0.00
17) Chrysene-d12	21.433	240	12708	0.400	ng/ul	0.00
23) Perylene-d12	23.824	264	10596	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	5221	1.341	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.250	152	25297	1.476	ng/ul	0.00
18) Fluoranthene-d10	19.276	212	53178	1.571	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	5575	1.380	ng/ul#	85
5) Naphthalene	10.711	128	46266	1.608	ng/ul	99
7) 2-Methylnaphthalene	12.322	142	31137	1.593	ng/ul	99
8) 1-Methylnaphthalene	12.542	142	30776	1.612	ng/ul	100
10) Acenaphthylene	14.219	152	37966	1.611	ng/ul	99
11) Acenaphthene	14.562	153	31179	1.625	ng/ul	99
12) Fluorene	15.547	166	36642	1.641	ng/ul	99
14) Pentachlorophenol	16.896	266	3387	0.955	ng/ul	99
15) Phenanthrene	17.289	178	58023	1.606	ng/ul	98
16) Anthracene	17.377	178	52438	1.645	ng/ul	97
19) Fluoranthene	19.304	202	67330	1.730	ng/ul	98
20) Pyrene	19.666	202	66162	1.684	ng/ul	99
21) Benzo(a)anthracene	21.419	228	60729	1.515	ng/ul	99
22) Chrysene	21.471	228	66068	1.580	ng/ul	100
24) Benzo(b)fluoranthene	23.093	252	63623	1.676	ng/ul	95
25) Benzo(k)fluoranthene	23.143	252	63064	1.676	ng/ul	96
26) Benzo(a)pyrene	23.716	252	53119	1.605	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	26.293	276	69068	1.591	ng/ul	99
28) Dibenzo(a,h)anthracene	26.314	278	57539	1.585	ng/ul	97
29) Benzo(g,h,i)perylene	27.051	276	58711	1.591	ng/ul	99

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

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