

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121621\
 Data File : BN018029.D
 Acq On : 15 Dec 2021 17:56
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV249

Quant Time: Dec 15 21:27:11 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN121621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 15 17:47:27 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.662	152	4288	0.400	ng/ul	0.00
4) Naphthalene-d8	10.458	136	12584	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.313	164	7868	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.070	188	15464	0.400	ng/ul	0.00
17) Chrysene-d12	21.272	240	16518	0.400	ng/ul #	0.00
23) Perylene-d12	23.543	264	14240	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.135	96	1798	0.404	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.064	152	5647	0.385	ng/ul	0.00
18) Fluoranthene-d10	19.105	212	19021	0.404	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.173	88	2351	0.450	ng/ul#	44
5) Naphthalene	10.508	128	13112	0.380	ng/ul	99
7) 2-Methylnaphthalene	12.141	142	7448	0.384	ng/ul	100
8) 1-Methylnaphthalene	12.350	142	8622	0.383	ng/ul	99
10) Acenaphthylene	14.031	152	12291	0.386	ng/ul	99
11) Acenaphthene	14.373	153	11046	0.390	ng/ul	97
12) Fluorene	15.372	166	12207	0.391	ng/ul	100
14) Pentachlorophenol	16.758	266	746	0.276	ng/ul	99
15) Phenanthrene	17.112	178	20846	0.389	ng/ul	99
16) Anthracene	17.205	178	16392	0.389	ng/ul	98
19) Fluoranthene	19.133	202	26663	0.401	ng/ul	99
20) Pyrene	19.495	202	30129	0.404	ng/ul	94
21) Benzo(a)anthracene	21.257	228	18177	0.370	ng/ul	99
22) Chrysene	21.304	228	27898	0.396	ng/ul	99
24) Benzo(b)fluoranthene	22.856	252	22463	0.395	ng/ul	85
25) Benzo(k)fluoranthene	22.903	252	29001	0.382	ng/ul#	82
26) Benzo(a)pyrene	23.446	252	21004	0.382	ng/ul#	82
27) Indeno(1,2,3-cd)pyrene	25.880	276	22822	0.369	ng/ul#	91
28) Dibenzo(a,h)anthracene	25.910	278	16503	0.365	ng/ul#	86
29) Benzo(g,h,i)perylene	26.581	276	20988	0.396	ng/ul#	88

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

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