

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090221\
 Data File : BN016151.D
 Acq On : 03 Sep 2021 00:11
 Operator : CG/JU
 Sample : PB138633BS
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS633

Quant Time: Sep 03 06:10:59 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 : Thu Sep 02 13:32:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.867	152	2545	0.400	ng/ul	0.00
4) Naphthalene-d8	10.666	136	11006	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.507	164	6625	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.255	188	13323	0.400	ng/ul	0.00
17) Chrysene-d12	21.441	240	13388	0.400	ng/ul	0.00
23) Perylene-d12	23.838	264	13321	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	2062	0.696	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	6812	0.439	ng/ul	0.00
18) Fluoranthene-d10	19.286	212	16324	0.454	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	5362	1.656	ng/ul#	88
5) Naphthalene	10.716	128	11103	0.379	ng/ul	99
7) 2-Methylnaphthalene	12.333	142	7798	0.408	ng/ul	100
8) 1-Methylnaphthalene	12.553	142	7290	0.382	ng/ul	99
10) Acenaphthylene	14.225	152	11238	0.448	ng/ul	100
11) Acenaphthene	14.567	153	7947	0.379	ng/ul	99
12) Fluorene	15.557	166	9092	0.379	ng/ul	99
14) Pentachlorophenol	16.905	266	2148	0.995	ng/ul	99
15) Phenanthrene	17.293	178	14717	0.383	ng/ul	99
16) Anthracene	17.386	178	14193	0.428	ng/ul	99
19) Fluoranthene	19.314	202	17933	0.402	ng/ul	99
20) Pyrene	19.676	202	18622	0.412	ng/ul	96
21) Benzo(a)anthracene	21.426	228	19128	0.467	ng/ul	98
22) Chrysene	21.479	228	19035	0.412	ng/ul	98
24) Benzo(b)fluoranthene	23.104	252	19344	0.386	ng/ul	91
25) Benzo(k)fluoranthene	23.154	252	19044	0.378	ng/ul#	94
26) Benzo(a)pyrene	23.730	252	17320	0.407	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	26.316	276	22105	0.402	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.336	278	18415	0.407	ng/ul#	94
29) Benzo(g,h,i)perylene	27.081	276	18703	0.392	ng/ul	97

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

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