

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN010322\
 Data File : BN018322.D
 Acq On : 03 Jan 2022 23:51
 Operator : CG/JU
 Sample : M5090-04
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BG3C6

Quant Time: Jan 04 01:11:52 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 28 00:34:45 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.607	152	1715	0.400	ng/ul	0.00
4) Naphthalene-d8	10.383	136	6445	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.249	164	5092	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	16.987	188	13451	0.400	ng/ul	0.00
17) Chrysene-d12	21.183	240	18540	0.400	ng/ul	0.00
23) Perylene-d12	23.404	264	27553	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.067	96	6829	4.265	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.994	152	3239	0.289	ng/ul	0.00
18) Fluoranthene-d10	19.023	212	18290	0.333	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.432	128	689	0.036	ng/ul#	72
7) 2-Methylnaphthalene	12.066	142	384	0.027	ng/ul	86
12) Fluorene	15.300	166	511	0.023	ng/ul#	91
14) Pentachlorophenol	16.658	266	157	0.029	ng/ul#	84
15) Phenanthrene	17.029	178	1608	0.036	ng/ul#	82
16) Anthracene	17.118	178	1127	0.028	ng/ul#	73
19) Fluoranthene	19.055	202	2331	0.032	ng/ul#	89
20) Pyrene	19.413	202	2169	0.028	ng/ul#	78
21) Benzo(a)anthracene	21.171	228	2023	0.029	ng/ul#	84
22) Chrysene	21.221	228	2348	0.031	ng/ul#	85
24) Benzo(b)fluoranthene	22.734	252	2702	0.023	ng/ul#	42
25) Benzo(k)fluoranthene	22.778	252	2677	0.020	ng/ul#	30
26) Benzo(a)pyrene	23.302	252	2439	0.022	ng/ul#	39
27) Indeno(1,2,3-cd)pyrene	25.657	276	3454	0.025	ng/ul#	89
28) Dibenzo(a,h)anthracene	25.681	278	2471	0.023	ng/ul#	42
29) Benzo(g,h,i)perylene	26.351	276	3089	0.026	ng/ul#	78

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

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