

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082322\
 Data File : BN021326.D
 Acq On : 23 Aug 2022 13:01
 Operator : CG/JU
 Sample : N4181-08
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X3C33

Quant Time: Aug 23 13:28:24 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN082222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 22 18:06:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.093	152	2789	0.400	ng/u1	0.00
4) Naphthalene-d8	10.927	136	8861	0.400	ng/u1	# 0.00
9) Acenaphthene-d10	14.749	164	4694	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.502	188	10107	0.400	ng/u1	0.00
17) Chrysene-d12	21.691	240	9493	0.400	ng/u1	# 0.00
23) Perylene-d12	24.225	264	8221	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.385	96	1394	0.387	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.511	152	3484	0.260	ng/u1	0.00
18) Fluoranthene-d10	19.524	212	10363	0.354	ng/u1	0.00
Target Compounds						
						Qvalue
5) Naphthalene	10.977	128	22995	0.820	ng/u1	97
10) Acenaphthylene	14.472	152	8547	0.353	ng/u1	99
11) Acenaphthene	14.814	153	8626	0.465	ng/u1	99
12) Fluorene	15.799	166	7711	0.372	ng/u1	99
14) Pentachlorophenol	17.143	266	498	0.446	ng/u1	98
15) Phenanthrene	17.540	178	16230	0.433	ng/u1	100
16) Anthracene	17.633	178	12303	0.410	ng/u1	99
19) Fluoranthene	19.557	202	21587	0.480	ng/u1	99
20) Pyrene	19.919	202	28123	0.624	ng/u1	98
21) Benzo(a)anthracene	21.673	228	19202	0.584	ng/u1	99
24) Benzo(b)fluoranthene	23.445	252	21148	0.599	ng/u1	97
27) Indeno(1,2,3-cd)pyrene	26.904	276	30977	0.793	ng/u1#	98
28) Dibenzo(a,h)anthracene	26.924	278	24345	0.758	ng/u1	94
29) Benzo(g,h,i)perylene	27.726	276	12178	0.314	ng/u1	97

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

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