

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030923\
 Data File : BN024195.D
 Acq On : 09 Mar 2023 11:02
 Operator : CG/JU
 Sample : 01746-19
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DCAH0

Quant Time: Mar 10 04:47:52 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN030923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 10 03:43:59 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.051	152	7640	0.400	ng/ul	0.00
4) Naphthalene-d8	10.872	136	21739	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.689	164	10846	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.434	188	22091	0.400	ng/ul	0.00
17) Chrysene-d12	21.627	240	14843	0.400	ng/ul	0.02
23) Perylene-d12	24.155	264	11655	0.400	ng/ul	# 0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.406	96	4826	0.574	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.451	152	8477	0.328	ng/ul	0.00
18) Fluoranthene-d10	19.459	212	15004	0.380	ng/ul	0.00
Target Compounds						
						Qvalue
5) Naphthalene	10.922	128	18085	0.333	ng/ul	100
7) 2-Methylnaphthalene	12.522	142	20106	0.611	ng/ul	100
10) Acenaphthylene	14.411	152	11793	0.261	ng/ul	99
12) Fluorene	15.735	166	20250	0.520	ng/ul	98
14) Pentachlorophenol	17.076	266	3683	0.600	ng/ul#	73
15) Phenanthrene	17.472	178	21534	0.340	ng/ul	100
16) Anthracene	17.565	178	27972	0.504	ng/ul	100
19) Fluoranthene	19.487	202	19743	0.334	ng/ul	98
20) Pyrene	19.850	202	25486	0.425	ng/ul	97
21) Benzo(a)anthracene	21.609	228	22075	0.456	ng/ul	100
22) Chrysene	21.665	228	25794	0.504	ng/ul	100
24) Benzo(b)fluoranthene	23.375	252	36468	0.819	ng/ul	99
25) Benzo(k)fluoranthene	23.427	252	17295	0.374	ng/ul	97
26) Benzo(a)pyrene	24.033	252	13892	0.363	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.851	278	22699	0.692	ng/ul	98
29) Benzo(g,h,i)perylene	27.643	276	13308	0.356	ng/ul	99

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

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