

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022824\
 Data File : BN030152.D
 Acq On : 29 Feb 2024 05:15
 Operator : MA/JU
 Sample : P1494-15MS
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BH3W6MS

Quant Time: Feb 29 05:54:41 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN020624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 23:49:29 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.871	152	3638	0.400	ng/ul	0.02
4) Naphthalene-d8	10.672	136	10141	0.400	ng/ul	0.02
9) Acenaphthene-d10	14.516	164	4458	0.400	ng/ul	0.02
13) Phenanthrene-d10	17.268	188	7366	0.400	ng/ul	0.00
17) Chrysene-d12	21.470	240	4020	0.400	ng/ul	0.00
23) Perylene-d12	23.841	264	3751	0.400	ng/ul	# 0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.331	96	1634	0.360	ng/ul	0.01
6) 2-Methylnaphthalene-d10	12.267	152	5522	0.428	ng/ul	0.02
18) Fluoranthene-d10	19.300	212	6876	0.520	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.361	88	4624	0.936	ng/ul	90
5) Naphthalene	10.716	128	11693	0.410	ng/ul	99
7) 2-Methylnaphthalene	12.338	142	6434	0.393	ng/ul	100
8) 1-Methylnaphthalene	12.558	142	6703	0.402	ng/ul	99
10) Acenaphthylene	14.234	152	8391	0.386	ng/ul	96
11) Acenaphthene	14.576	153	6913	0.418	ng/ul	98
12) Fluorene	15.571	166	6979	0.404	ng/ul	99
14) Pentachlorophenol	16.918	266	1843	1.122	ng/ul#	50
15) Phenanthrene	17.310	178	10823	0.470	ng/ul	98
16) Anthracene	17.407	178	8739	0.451	ng/ul	98
19) Fluoranthene	19.332	202	8515	0.461	ng/ul	95
20) Pyrene	19.695	202	8994	0.475	ng/ul#	92
21) Benzo(a)anthracene	21.453	228	6581	0.454	ng/ul	98
22) Chrysene	21.505	228	7477	0.471	ng/ul	98
24) Benzo(b)fluoranthene	23.119	252	7195	0.510	ng/ul	94
25) Benzo(k)fluoranthene	23.168	252	8117	0.519	ng/ul	95
26) Benzo(a)pyrene	23.738	252	6683	0.504	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	26.309	276	7702	0.478	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.349	278	5949	0.471	ng/ul#	89
29) Benzo(g,h,i)perylene	27.050	276	7186	0.491	ng/ul	94

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

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