

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032922\
 Data File : BM034405.D
 Acq On : 29 Mar 2022 13:52
 Operator : CG/JU
 Sample : N2079-20MS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW607MS

Quant Time: Mar 30 02:29:59 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM032822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 28 18:27:40 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.933	152	2300	0.400	ng/ul	0.00
4) Naphthalene-d8	10.741	136	6695	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.573	164	3792	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.316	188	6829	0.400	ng/ul	0.00
17) Chrysene-d12	21.495	240	5723	0.400	ng/ul	# 0.00
23) Perylene-d12	23.903	264	6360	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	8712	2.546	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.335	152	3234	0.358	ng/ul	0.00
18) Fluoranthene-d10	19.343	212	8182	0.458	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.343	88	4365	1.155	ng/ul#	79
5) Naphthalene	10.790	128	46453	2.396	ng/ul#	88
7) 2-Methylnaphthalene	12.407	142	7781	0.656	ng/ul	99
8) 1-Methylnaphthalene	12.627	142	6995	0.591	ng/ul	99
10) Acenaphthylene	14.296	152	6444	0.353	ng/ul	92
11) Acenaphthene	14.634	153	4598	0.331	ng/ul	96
12) Fluorene	15.624	166	5066	0.342	ng/ul	96
14) Pentachlorophenol	16.978	266	2414	0.878	ng/ul	98
15) Phenanthrene	17.358	178	7877	0.361	ng/ul	97
16) Anthracene	17.455	178	6546	0.342	ng/ul	95
19) Fluoranthene	19.371	202	10179	0.395	ng/ul	99
20) Pyrene	19.733	202	11144	0.391	ng/ul	97
21) Benzo(a)anthracene	21.477	228	6691	0.374	ng/ul	99
22) Chrysene	21.530	228	8196	0.400	ng/ul	98
24) Benzo(b)fluoranthene	23.167	252	8267	0.376	ng/ul	84
25) Benzo(k)fluoranthene	23.213	252	7753	0.364	ng/ul#	85
26) Benzo(a)pyrene	23.798	252	9400	0.412	ng/ul#	82
27) Indeno(1,2,3-cd)pyrene	26.418	276	10870	0.366	ng/ul#	90
28) Dibenzo(a,h)anthracene	26.441	278	8386	0.376	ng/ul#	83
29) Benzo(g,h,i)perylene	27.186	276	11134	0.388	ng/ul#	95

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

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