

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071723\
 Data File : BN026544.D
 Acq On : 17 Jul 2023 19:38
 Operator : MA/JU
 Sample : PB154018BS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS018

Quant Time: Jul 18 04:21:56 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN071223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 17 17:14:21 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.967	152	1714	0.400	ng/ul	0.02
4) Naphthalene-d8	10.765	136	5076	0.400	ng/ul	0.01
9) Acenaphthene-d10	14.584	164	3497	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.334	188	6948	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.520	240	5555	0.400	ng/ul	# 0.00
23) Perylene-d12	23.932	264	5409	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.297	96	2022	0.896	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.359	152	2794	0.405	ng/ul	0.01
18) Fluoranthene-d10	19.359	212	8133	0.422	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.335	88	4725	2.212	ng/ul	89
5) Naphthalene	10.814	128	4498	0.323	ng/ul	95
7) 2-Methylnaphthalene	12.431	142	2676	0.338	ng/ul	99
8) 1-Methylnaphthalene	12.640	142	3227	0.381	ng/ul	97
10) Acenaphthylene	14.311	152	5413	0.316	ng/ul	99
11) Acenaphthene	14.644	153	4135	0.317	ng/ul	98
12) Fluorene	15.639	166	4629	0.338	ng/ul	99
14) Pentachlorophenol	17.005	266	1114	0.636	ng/ul	97
15) Phenanthrene	17.376	178	6973	0.315	ng/ul	99
16) Anthracene	17.473	178	6156	0.321	ng/ul	99
19) Fluoranthene	19.391	202	8797	0.338	ng/ul#	93
20) Pyrene	19.754	202	9523	0.337	ng/ul#	94
21) Benzo(a)anthracene	21.505	228	6564	0.322	ng/ul	99
22) Chrysene	21.555	228	7683	0.331	ng/ul	99
24) Benzo(b)fluoranthene	23.195	252	6685	0.317	ng/ul	95
25) Benzo(k)fluoranthene	23.245	252	7727	0.350	ng/ul#	94
26) Benzo(a)pyrene	23.826	252	6359	0.339	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	26.453	276	8950	0.364	ng/ul#	75
28) Dibenzo(a,h)anthracene	26.464	278	7026	0.372	ng/ul#	78
29) Benzo(g,h,i)perylene	27.228	276	7858	0.362	ng/ul#	88

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

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