

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070324\
 Data File : BN032435.D
 Acq On : 03 Jul 2024 13:45
 Operator : MA/JU
 Sample : P2964-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4DN9

Quant Time: Jul 03 14:10:58 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN062724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 03 03:26:42 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	6736	0.400	ng/ul	0.00
4) Naphthalene-d8	10.376	136	21269	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.248	164	12781	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.003	188	25203	0.400	ng/ul	0.00
17) Chrysene-d12	21.211	240	18078	0.400	ng/ul	# 0.00
23) Perylene-d12	23.433	264	20340	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	2818	0.376	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.977	152	7344	0.230	ng/ul	0.00
18) Fluoranthene-d10	19.045	212	19221	0.353	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.426	128	502939	8.935	ng/ul	98
10) Acenaphthylene	13.966	152	225617	4.244	ng/ul	99
11) Acenaphthene	14.309	153	214556	5.201	ng/ul	98
12) Fluorene	15.308	166	76034	1.547	ng/ul	98
15) Phenanthrene	17.046	178	842294	12.190	ng/ul	99
16) Anthracene	17.138	178	679567	11.026	ng/ul	99
19) Fluoranthene	19.073	202	182868	2.656	ng/ul	98
20) Pyrene	19.440	202	275078	3.824	ng/ul	98
21) Benzo(a)anthracene	21.197	228	101199	1.551	ng/ul	100
22) Chrysene	21.249	228	535305	7.475	ng/ul	100
24) Benzo(b)fluoranthene	22.763	252	266330	3.014	ng/ul	94
25) Benzo(k)fluoranthene	22.807	252	67344	0.784	ng/ul#	88
26) Benzo(a)pyrene	23.333	252	343077	5.580	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	25.666	276	223604	2.323	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.673	278	194786	2.723	ng/ul	97
29) Benzo(g,h,i)perylene	26.350	276	550073	7.173	ng/ul	98

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

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