

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022324\
 Data File : BN029960.D
 Acq On : 23 Feb 2024 17:16
 Operator : MA/JU
 Sample : P1506-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BP-VPB-195-GW-840-842

Quant Time: Feb 23 17:43:30 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN012924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 16 23:52:05 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.862	152	4521	0.400	ng	0.00
7) Naphthalene-d8	10.648	136	11554	0.400	ng	#-0.02
13) Acenaphthene-d10	14.500	164	5595	0.400	ng	-0.01
19) Phenanthrene-d10	17.255	188	10242	0.400	ng	0.00
29) Chrysene-d12	21.456	240	9408	0.400	ng	# 0.00
35) Perylene-d12	23.820	264	9120	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.435	112	2192	0.202	ng	0.00
5) Phenol-d6	7.031	99	1980	0.158	ng	0.00
8) Nitrobenzene-d5	9.025	82	4095	0.426	ng	-0.01
11) 2-Methylnaphthalene-d10	12.246	152	5644	0.359	ng	-0.01
14) 2,4,6-Tribromophenol	16.002	330	922	0.514	ng	0.00
15) 2-Fluorobiphenyl	13.121	172	9114	0.379	ng	-0.01
27) Fluoranthene-d10	19.289	212	11091	0.447	ng	0.00
31) Terphenyl-d14	19.898	244	9132	0.391	ng	0.00
Target Compounds						
						Qvalue
33) Chrysene	21.492	228	813	0.022	ng	# 81
34) Bis(2-ethylhexyl)phtha...	21.384	149	37207	1.699	ng	98
37) Benzo(b)fluoranthene	23.104	252	1097	0.034	ng	# 1
38) Benzo(k)fluoranthene	23.151	252	930	0.027	ng	# 1
41) Benzo(g,h,i)perylene	27.025	276	735	0.021	ng	# 36

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

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